

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062798.D
 Acq On : 13 Sep 2024 21:36
 Operator : JU/RC
 Sample : P3898-20
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC59

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062798.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.937	478	485	494	rVB2	5716566	10774220	5.42%	1.038%
2	6.407	558	565	572	rBV4	1657704	3490366	1.75%	0.336%
3	6.478	572	577	582	rBV	2196007	3675992	1.85%	0.354%
4	6.560	582	591	594	rBV2	1750024	4273725	2.15%	0.412%
5	6.918	640	652	657	rBV3	2978074	9145267	4.60%	0.881%
6	6.977	657	662	665	rVV	4022304	6931678	3.48%	0.668%
7	7.012	665	668	673	rVV	3759814	6278648	3.16%	0.605%
8	7.077	673	679	685	rVV3	5030688	12186557	6.13%	1.174%
9	7.147	685	691	695	rVV	2677760	4581522	2.30%	0.441%
10	7.306	712	718	722	rVV	2850481	5417713	2.72%	0.522%
11	7.353	722	726	732	rVV	4240269	7852190	3.95%	0.757%
12	7.435	733	740	745	rVV4	1510314	3856711	1.94%	0.372%
13	7.576	754	764	773	rVV	16545431	36978630	18.59%	3.563%
14	7.753	782	794	798	rBV4	1138372	3047024	1.53%	0.294%
15	7.847	801	810	815	rBV3	1157841	3368251	1.69%	0.325%
16	7.911	815	821	825	rVB	4489192	7581302	3.81%	0.731%
17	8.011	825	838	842	rBV2	6912778	20087582	10.10%	1.936%
18	8.123	851	857	858	rBV2	1940981	3014073	1.52%	0.290%
19	8.152	858	862	865	rVB	3787090	5533713	2.78%	0.533%
20	8.364	891	898	901	rBV	4186203	7873074	3.96%	0.759%
21	8.464	910	915	921	rVB3	4255246	8061472	4.05%	0.777%
22	8.522	921	925	928	rBV	2714013	3561059	1.79%	0.343%
23	8.575	928	934	935	rBV2	4261751	6561300	3.30%	0.632%
24	8.699	948	955	958	rBV	4357927	7959446	4.00%	0.767%
25	8.728	958	960	969	rVB2	2919665	5081493	2.55%	0.490%
26	8.881	980	986	989	rBV	2617768	4490622	2.26%	0.433%
27	8.928	989	994	1001	rVB2	3188310	6721871	3.38%	0.648%
28	9.033	1001	1012	1021	rBV4	4955392	12764296	6.42%	1.230%
29	9.186	1021	1038	1044	rVB	14170188	38528202	19.37%	3.713%
30	9.280	1044	1054	1060	rBV7	3572992	11014342	5.54%	1.061%
31	9.357	1060	1067	1071	rBV5	2455838	5677825	2.85%	0.547%
32	9.509	1088	1093	1097	rVB2	1775014	2953892	1.48%	0.285%
33	9.562	1097	1102	1106	rBV2	2107635	3932786	1.98%	0.379%
34	9.615	1106	1111	1121	rVB3	2187822	5309405	2.67%	0.512%
35	9.715	1121	1128	1134	rVB2	2226962	4934838	2.48%	0.476%
36	9.803	1134	1143	1147	rBV4	2174023	5772985	2.90%	0.556%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
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37	9.856	1147	1152	1159	rBV4	3184832	8195447	4.12%	0.790%
38	10.003	1168	1177	1181	rVB	3793474	8734637	4.39%	0.842%
39	10.068	1181	1188	1192	rBV2	2634346	4327939	2.18%	0.417%
40	10.150	1193	1202	1205	rVV3	3355536	8378387	4.21%	0.807%
41	10.250	1213	1219	1223	rBV	2292541	4101696	2.06%	0.395%
42	10.408	1240	1246	1251	rBV2	2482263	4339829	2.18%	0.418%
43	10.726	1288	1300	1314	rVB3	12414055	45431484	22.84%	4.378%
44	10.843	1314	1320	1324	rBV2	7720584	13213668	6.64%	1.273%
45	10.884	1324	1327	1333	rVB	2874152	4169941	2.10%	0.402%
46	10.996	1341	1346	1357	rVB4	1581744	3829625	1.93%	0.369%
47	11.102	1357	1364	1375	rBV	8668602	16403285	8.25%	1.581%
48	11.231	1376	1386	1391	rBV	5548282	11120939	5.59%	1.072%
49	11.630	1446	1454	1468	rVB8	2508432	8693201	4.37%	0.838%
50	11.754	1468	1475	1479	rBV	6343430	11605485	5.83%	1.118%
51	11.818	1481	1486	1492	rVB	5801777	10943282	5.50%	1.055%
52	11.983	1505	1514	1522	rBV2	6380994	11928459	6.00%	1.149%
53	12.159	1534	1544	1546	rBV	8279681	17867708	8.98%	1.722%
54	12.183	1546	1548	1553	rVB	5141973	6582012	3.31%	0.634%
55	12.271	1553	1563	1566	rBV4	1982168	4559057	2.29%	0.439%
56	12.388	1576	1583	1588	rVB2	10173566	18753888	9.43%	1.807%
57	12.547	1604	1610	1614	rBV2	3833645	7093283	3.57%	0.684%
58	12.600	1615	1619	1626	rVB2	2388127	4784840	2.41%	0.461%
59	12.794	1644	1652	1657	rBV5	1694686	4768822	2.40%	0.460%
60	12.947	1674	1678	1686	rBV4	1811233	4523300	2.27%	0.436%
61	13.093	1698	1703	1706	rBV	2678097	4332681	2.18%	0.418%
62	13.305	1734	1739	1743	rVB	2707682	3847551	1.93%	0.371%
63	13.393	1746	1754	1759	rBV	8183150	15678961	7.88%	1.511%
64	13.522	1771	1776	1783	rVV	3365565	7578079	3.81%	0.730%
65	13.605	1783	1790	1796	rVB4	2789563	5999763	3.02%	0.578%
66	13.734	1805	1812	1816	rBV3	3066605	6098288	3.07%	0.588%
67	13.822	1821	1827	1831	rBV	3999294	6804010	3.42%	0.656%
68	13.881	1831	1837	1842	rBV3	3236226	7665263	3.85%	0.739%
69	14.039	1856	1864	1868	rBV2	3653569	7312684	3.68%	0.705%
70	14.075	1868	1870	1880	rVV4	1691962	4636181	2.33%	0.447%
71	14.180	1884	1888	1896	rVB4	4362388	9379486	4.72%	0.904%
72	14.286	1901	1906	1909	rBV3	1804684	3698064	1.86%	0.356%
73	14.474	1930	1938	1941	rBV	13783255	24327134	12.23%	2.344%
74	14.621	1959	1963	1971	rVB7	3440364	5893375	2.96%	0.568%
75	14.703	1971	1977	1982	rVB	8273657	12355211	6.21%	1.191%
76	14.762	1983	1987	1995	rBV4	2466846	5450936	2.74%	0.525%

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77	15.032	2025	2033	2036	rBV3	1479379	3735096	1.88%	0.360%
78	15.073	2037	2040	2044	rBV3	1601941	3064673	1.54%	0.295%
79	15.214	2055	2064	2067	rBV	13426762	28798196	14.48%	2.775%
80	15.326	2077	2083	2089	rBV2	8595908	14671998	7.38%	1.414%
81	15.414	2089	2098	2101	rVB	8400956	15114172	7.60%	1.456%
82	15.808	2162	2165	2169	rVB2	3117268	3914155	1.97%	0.377%
83	15.902	2177	2181	2187	rVB5	2093780	3560133	1.79%	0.343%
84	16.178	2223	2228	2238	rVB4	3448307	6385307	3.21%	0.615%
85	16.272	2238	2244	2246	rBV	8012454	13290961	6.68%	1.281%
86	16.607	2298	2301	2308	rBV3	1645430	2999745	1.51%	0.289%
87	17.071	2354	2380	2384	rBV3	26965806	198925126	100.00%	19.169%
88	17.600	2464	2470	2475	rBV4	2765610	5489378	2.76%	0.529%
89	17.800	2499	2504	2509	rVB	6819034	9492231	4.77%	0.915%
90	17.964	2528	2532	2538	rVB	4461672	5927846	2.98%	0.571%
91	18.487	2616	2621	2625	rBV	4971603	6415769	3.23%	0.618%
92	19.116	2721	2728	2730	rBV2	4518593	8038339	4.04%	0.775%
93	19.686	2821	2825	2829	rBV2	3486594	4019178	2.02%	0.387%
94	20.214	2911	2915	2918	rBV	2904025	3840185	1.93%	0.370%
95	20.708	2995	2999	3003	rBV	2555359	2901600	1.46%	0.280%
96	21.202	3079	3083	3096	rVB	5305619	7313452	3.68%	0.705%
97	21.724	3167	3172	3183	rVB	2075511	3296201	1.66%	0.318%
98	22.189	3245	3251	3258	rVB	2034093	3605544	1.81%	0.347%
99	22.306	3265	3271	3281	rVB3	2554284	5205169	2.62%	0.502%
100	24.603	3654	3662	3673	rVB3	881262	3031039	1.52%	0.292%

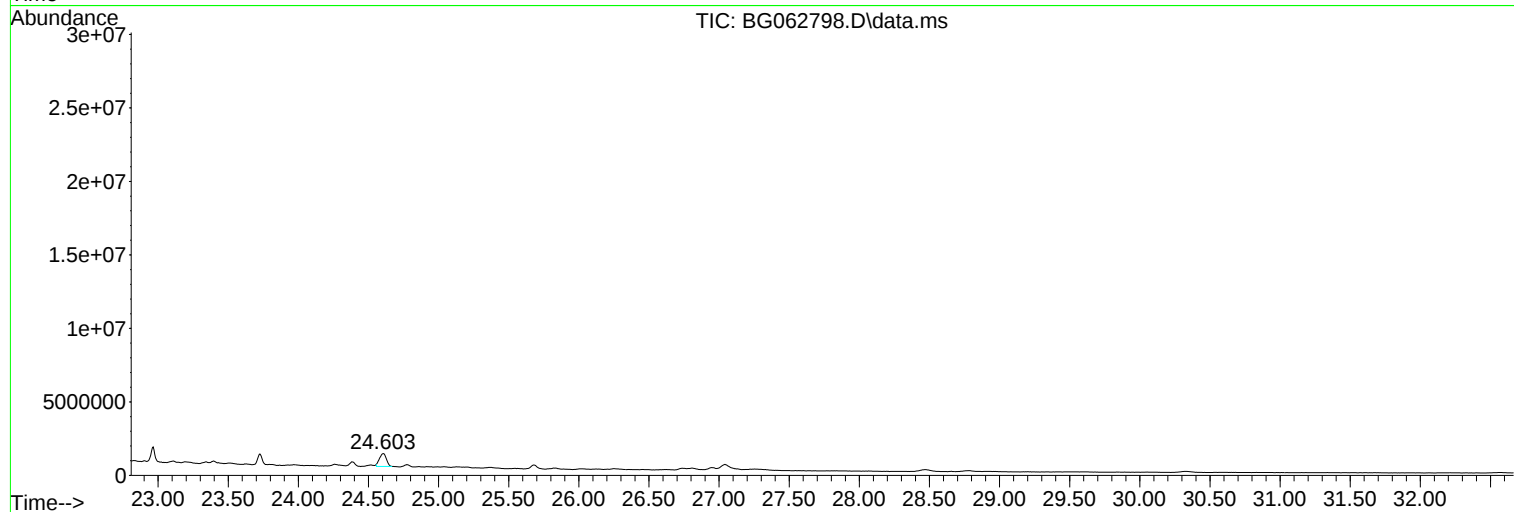
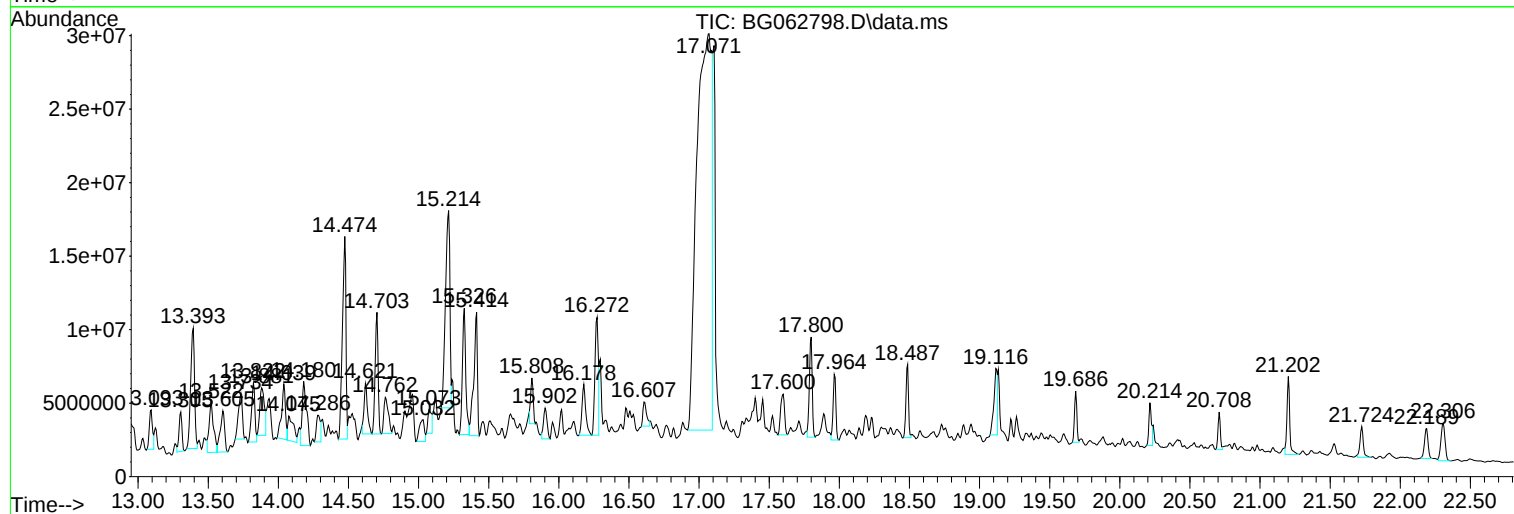
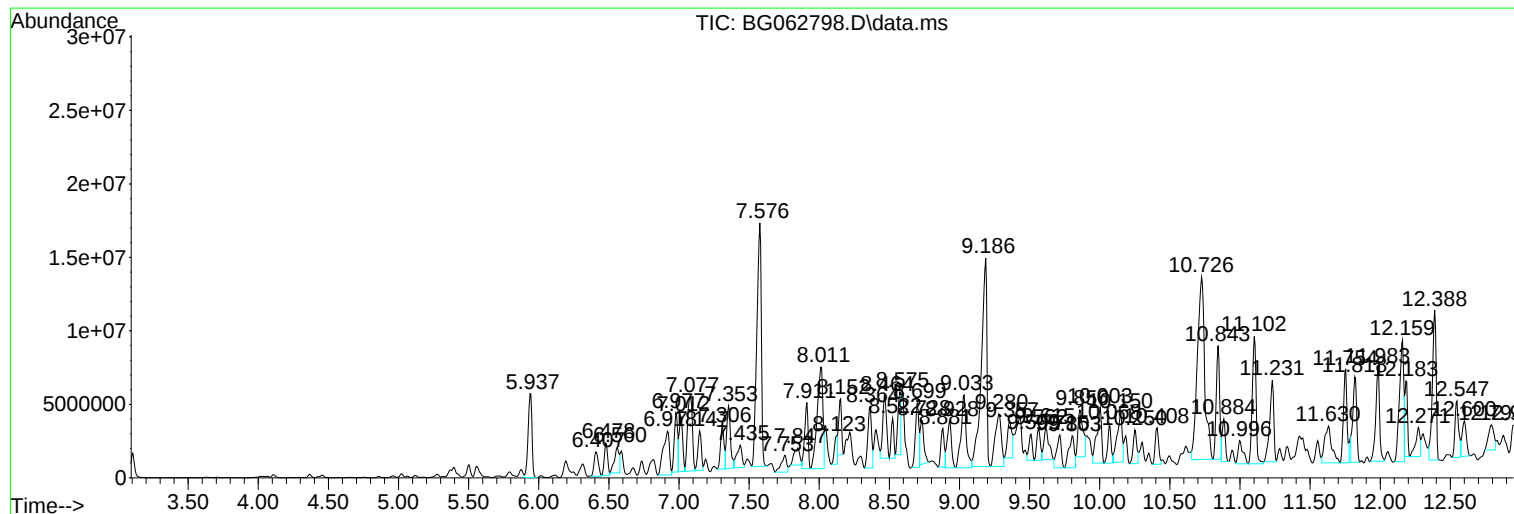
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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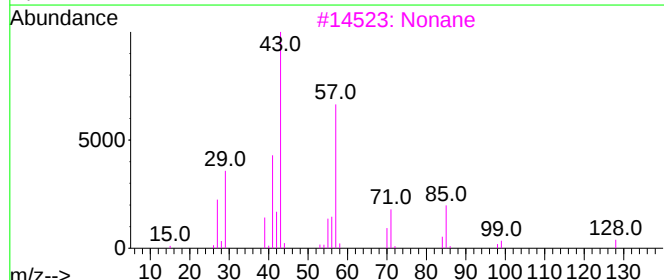
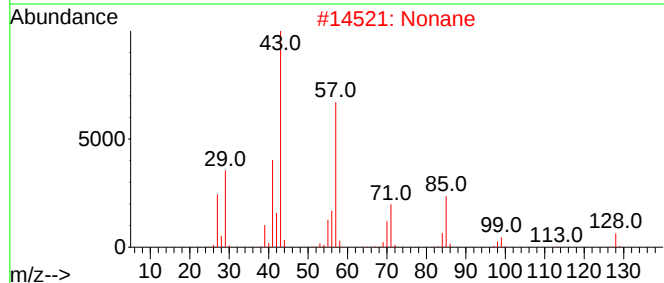
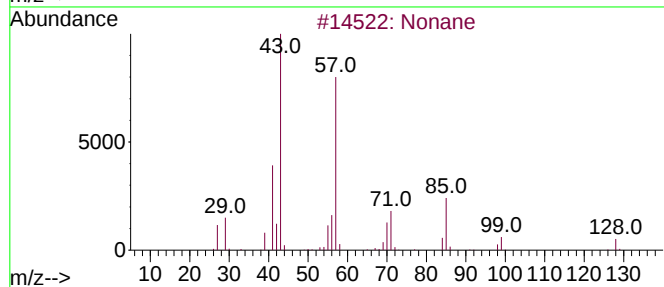
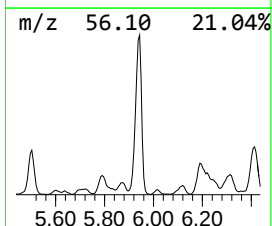
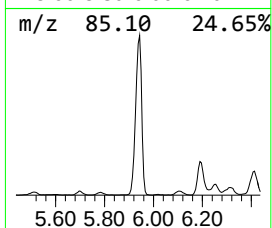
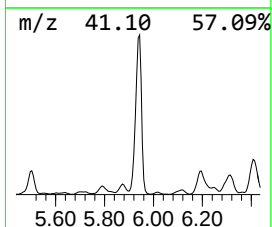
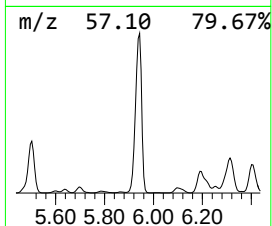
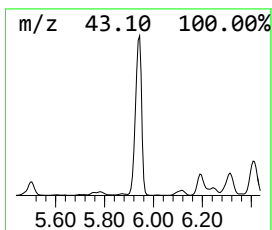
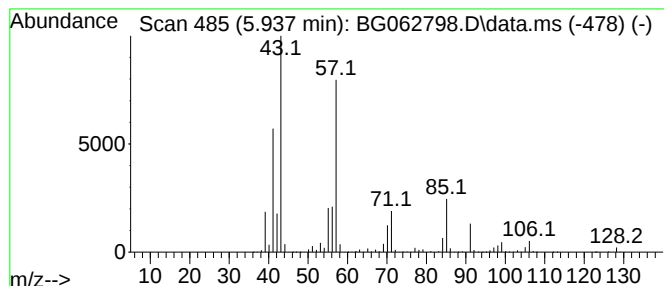
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.937	28.42 ng/ul	10774200	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	81
2			Nonane	128	C9H20	000111-84-2	80
3			Nonane	128	C9H20	000111-84-2	80
4			Carbonic acid, decyl prop-1-en-2...	242	C14H26O3	1000382-90-5	59
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59



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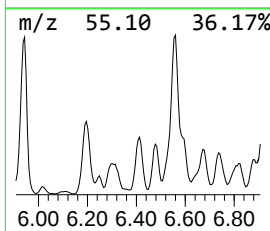
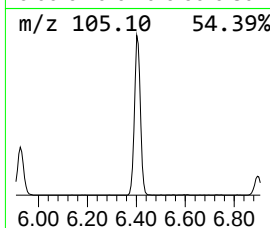
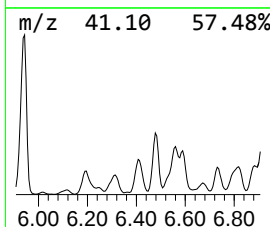
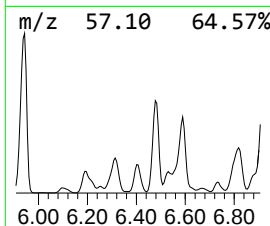
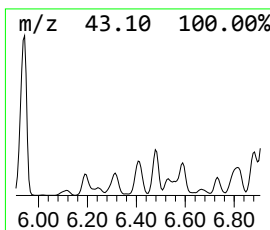
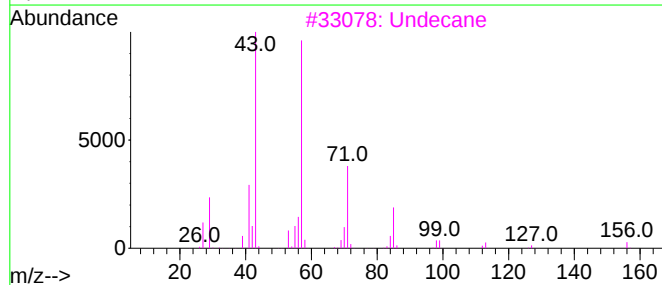
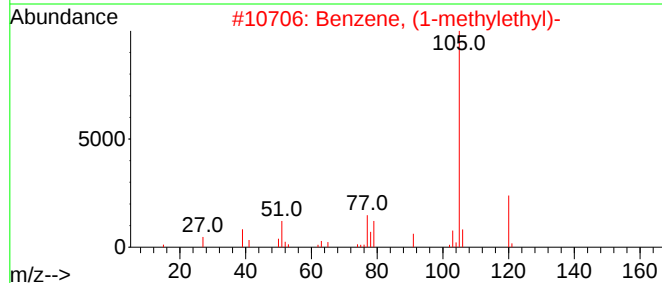
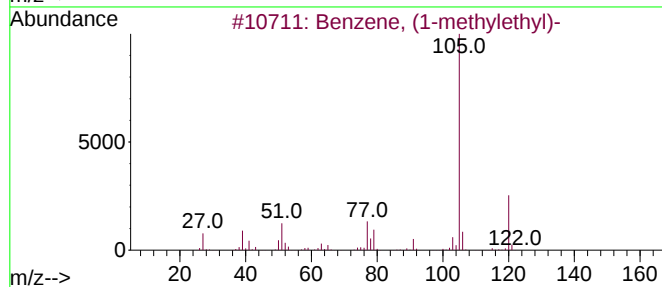
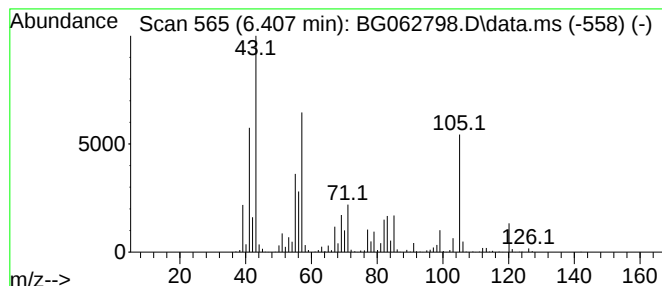
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, (1-methylethyl)- Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.407	9.21 ng/ul	3490370	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	53
2			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	53
3			Undecane	156	C11H24	001120-21-4	35
4			Octane	114	C8H18	000111-65-9	27
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	25



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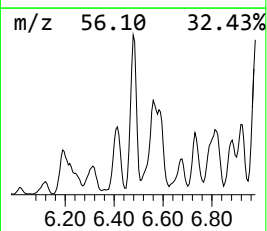
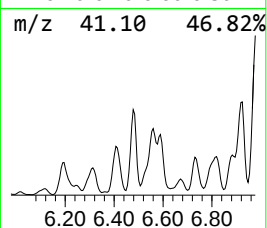
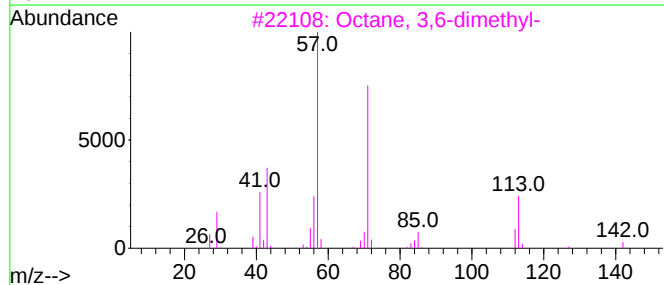
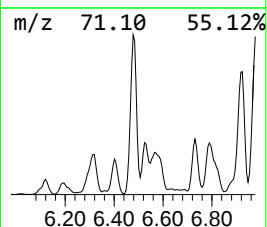
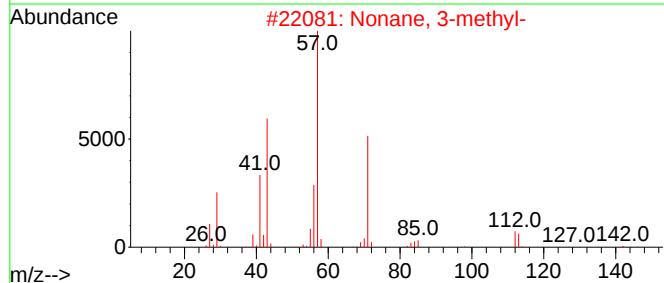
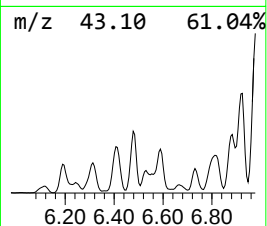
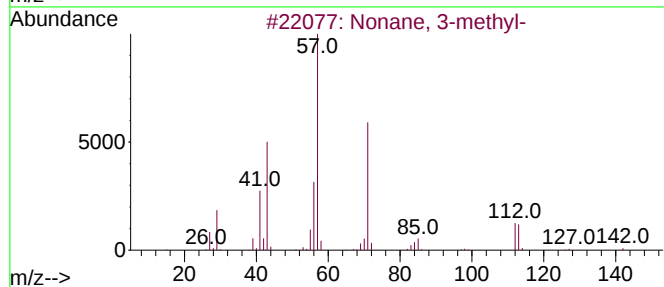
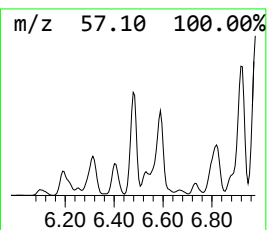
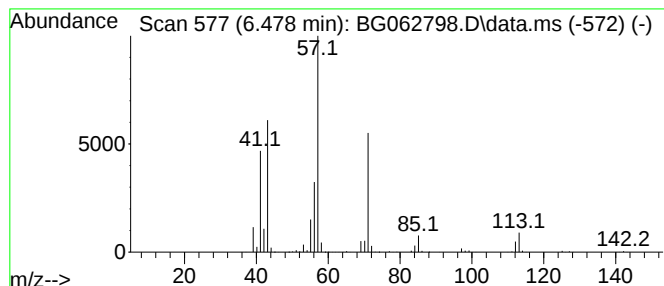
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.478	9.70 ng/ul	3675990	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	78
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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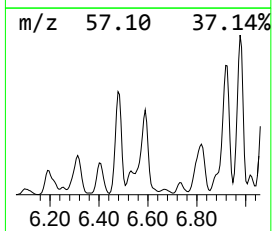
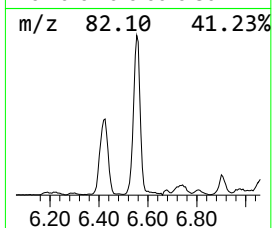
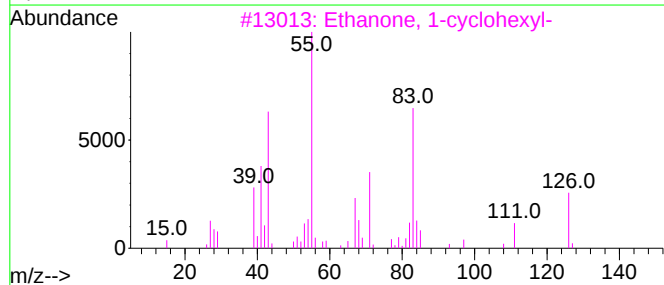
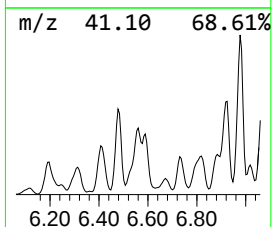
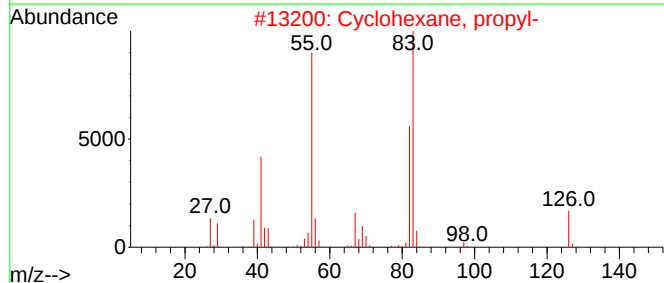
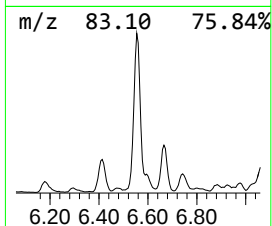
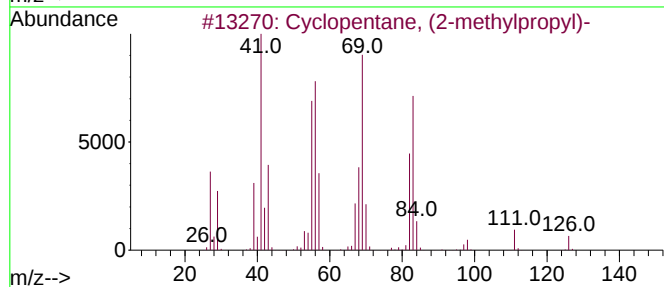
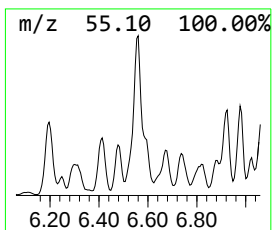
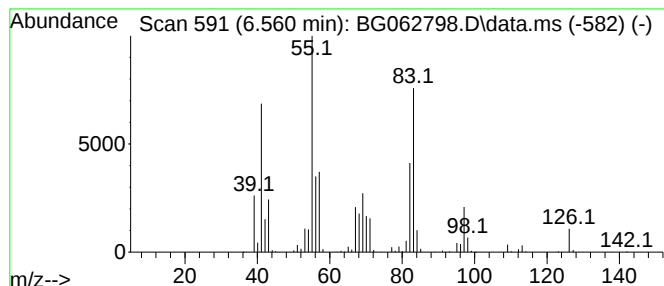
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic6.560 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.560	11.27 ng/ul	4273730	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, (2-methylpropyl)-	126	C9H18	003788-32-7	68
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
3			Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	50
4			1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	43
5			3,4-Dimethyl-2-hexene	112	C8H16	002213-37-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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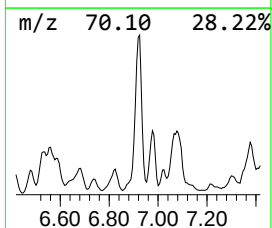
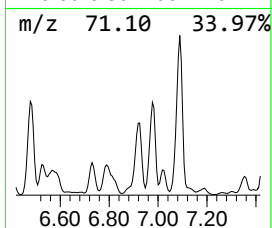
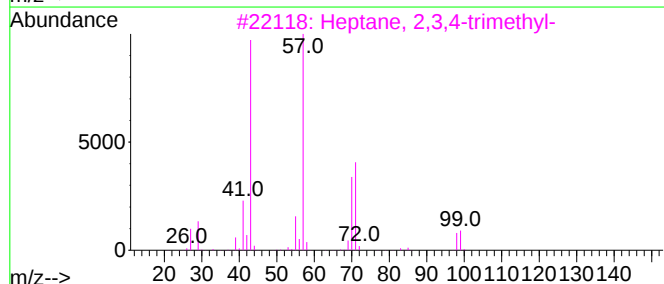
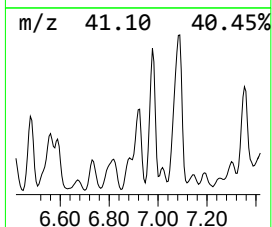
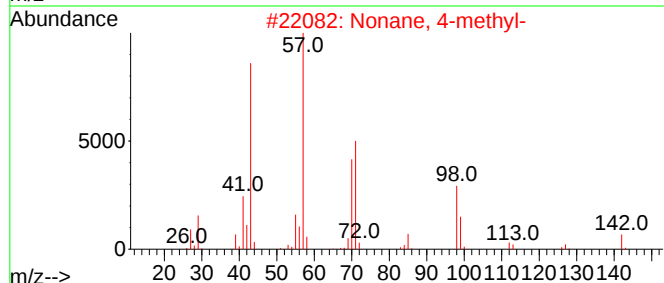
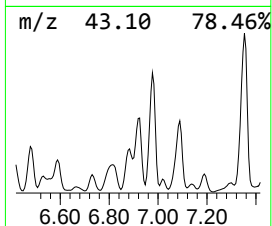
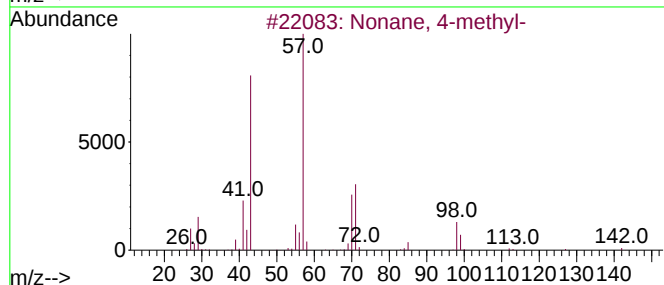
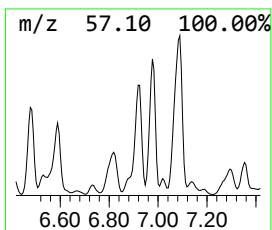
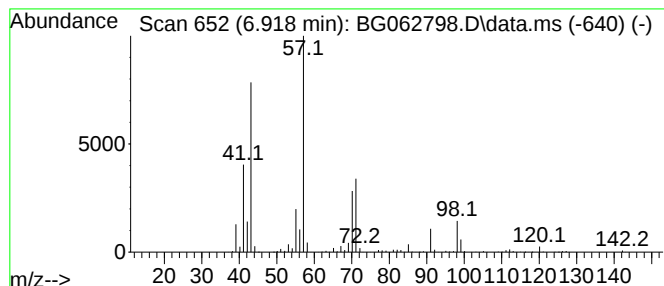
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.918	24.13 ng/ul	9145270	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	90
2			Nonane, 4-methyl-	142	C10H22	017301-94-9	78
3			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	59
4			Heptane, 2,3,5-trimethyl-	142	C10H22	020278-85-7	50
5			Octane, 2-bromo-	192	C8H17Br	000557-35-7	50



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Sample : P3898-20
Misc :
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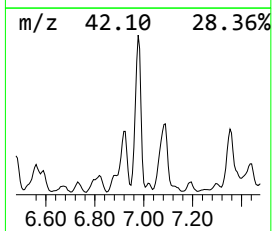
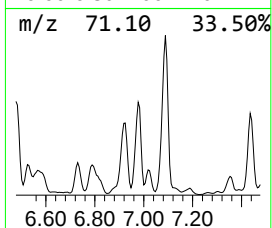
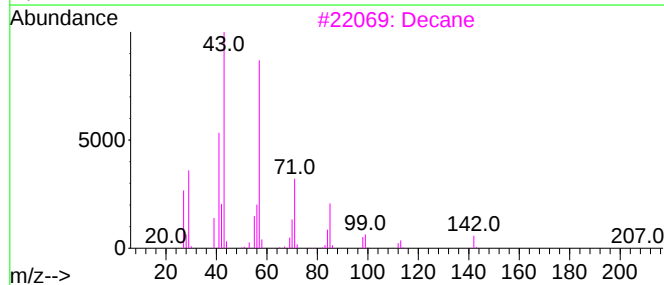
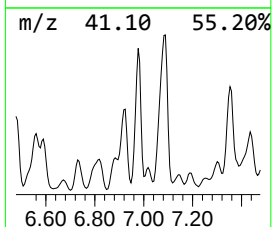
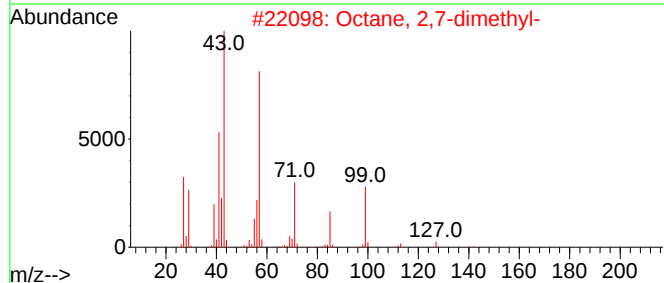
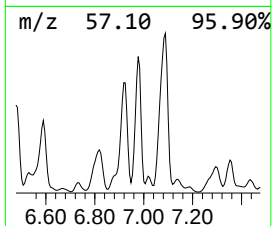
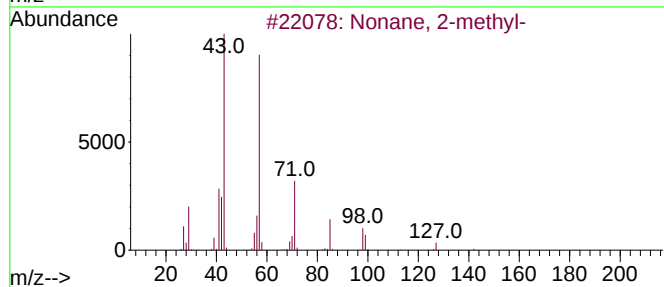
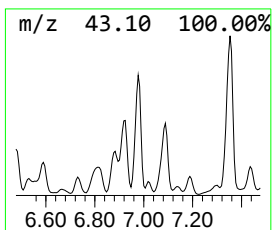
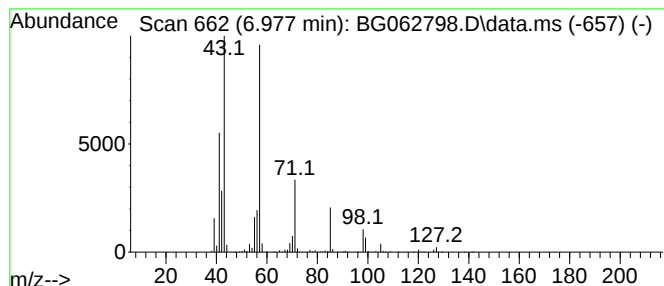
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.977	18.29 ng/ul	6931680	1,4-Dichlorobenzene-d4			7.917
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-		142	C10H22	000871-83-0	91
2	Octane, 2,7-dimethyl-		142	C10H22	001072-16-8	87
3	Decane		142	C10H22	000124-18-5	83
4	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	59
5	Octane, 2,7-dimethyl-		142	C10H22	001072-16-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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Acq On : 13 Sep 2024 21:36
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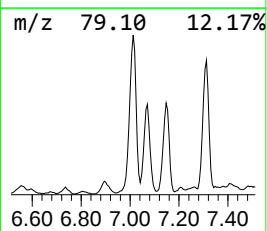
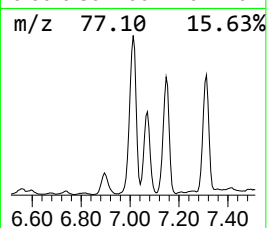
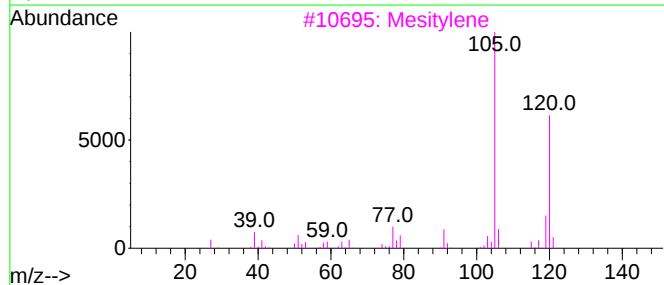
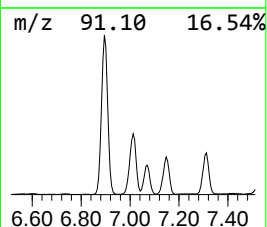
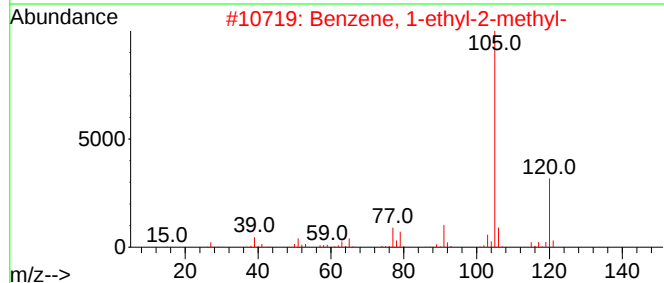
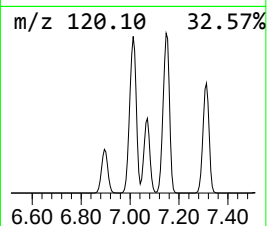
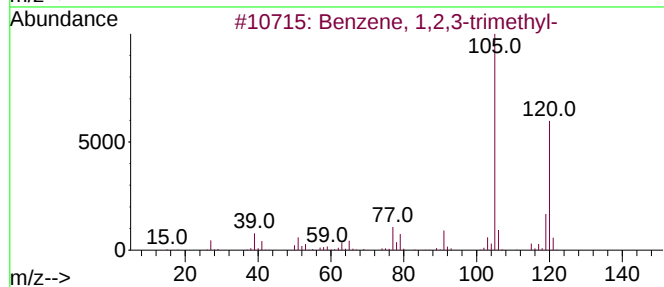
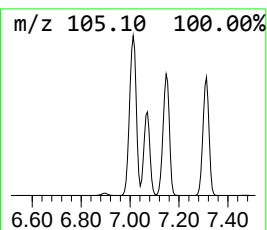
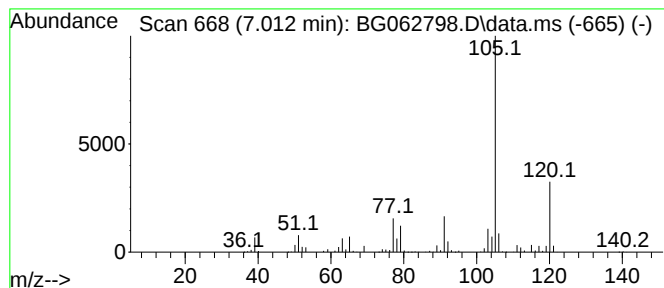
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.012	16.56 ng/ul	6278650	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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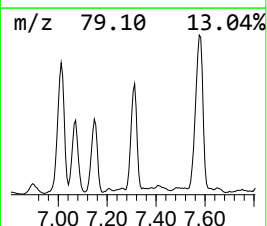
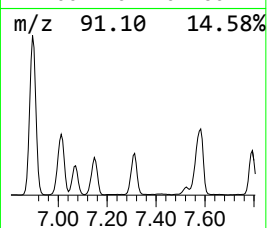
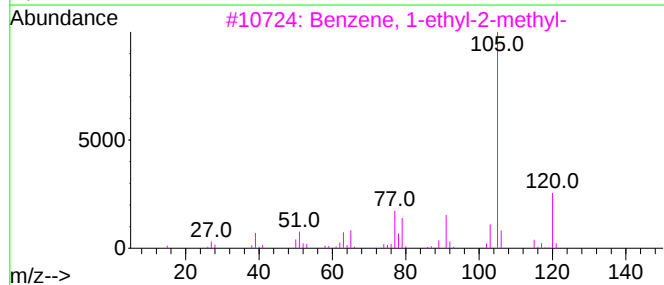
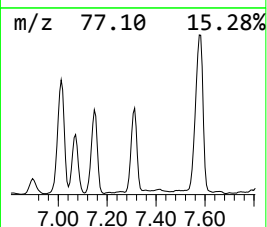
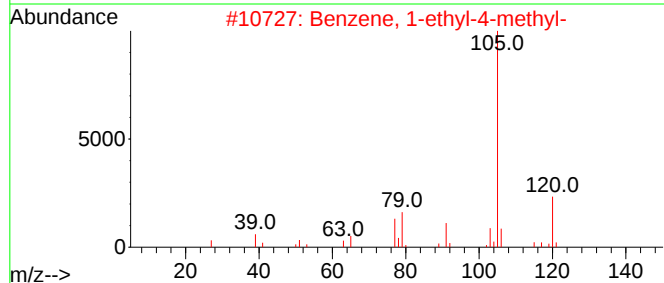
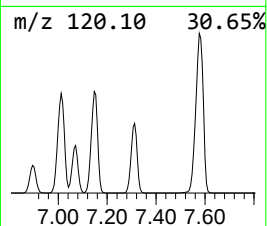
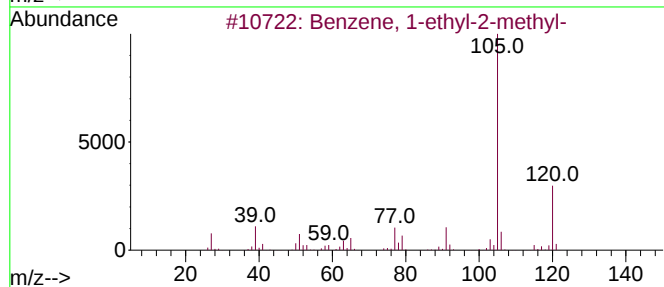
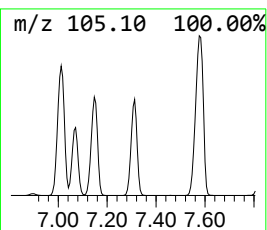
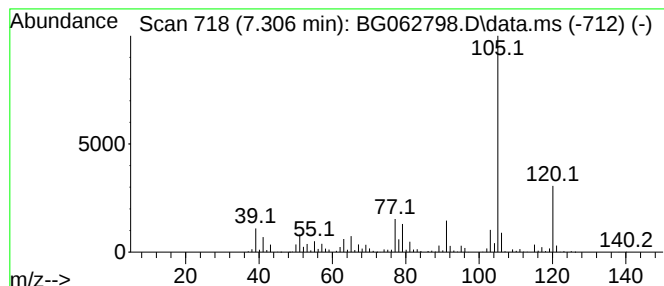
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.306	14.29 ng/ul	5417710	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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Misc :
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ClientSampleId :
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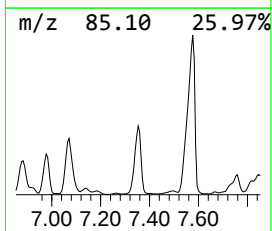
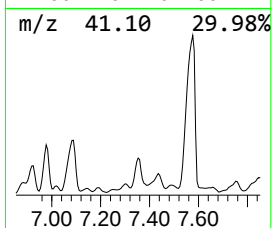
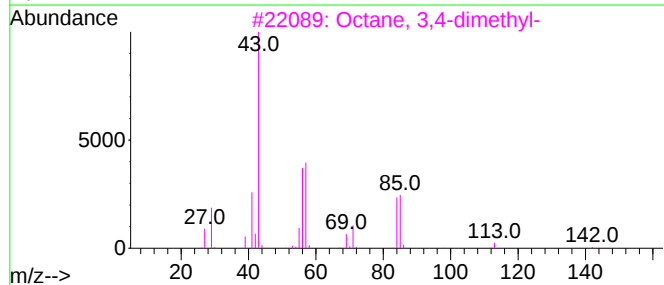
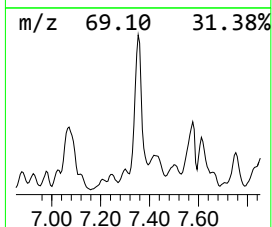
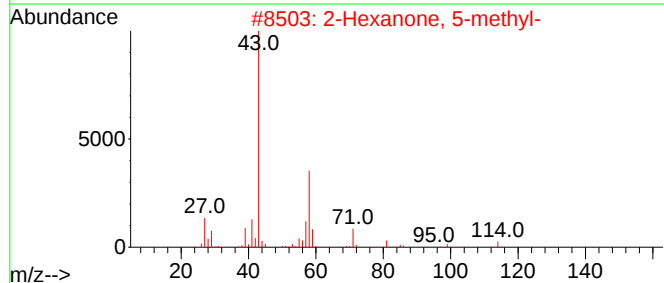
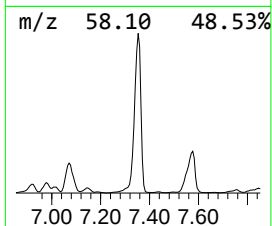
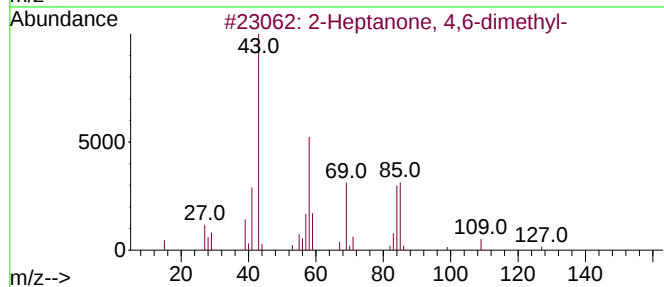
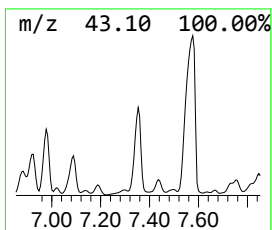
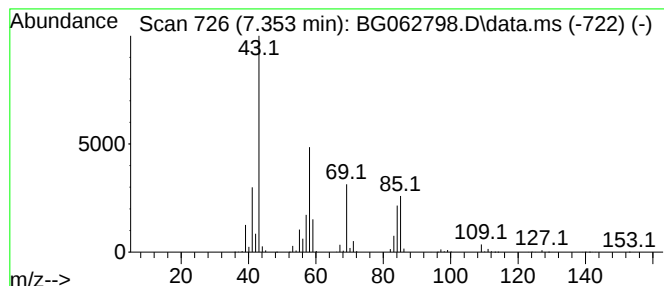
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 2-Heptanone, 4,6-dimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.353	20.71 ng/ul	7852190	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-			142	C9H18O	019549-80-5	83
2	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	47
3	Octane, 3,4-dimethyl-			142	C10H22	015869-92-8	43
4	Hexane, 3-ethyl-			114	C8H18	000619-99-8	43
5	2-Nonanone, 9-[(tetrahydro-2H-py...			242	C14H26O3	054699-41-1	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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ClientSampleId :
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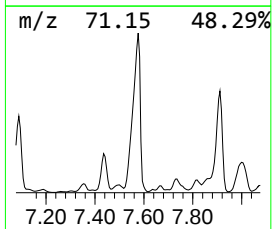
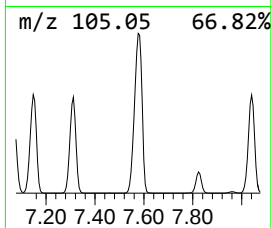
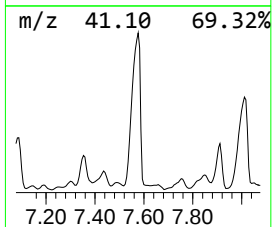
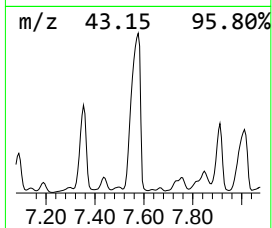
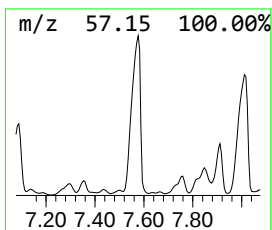
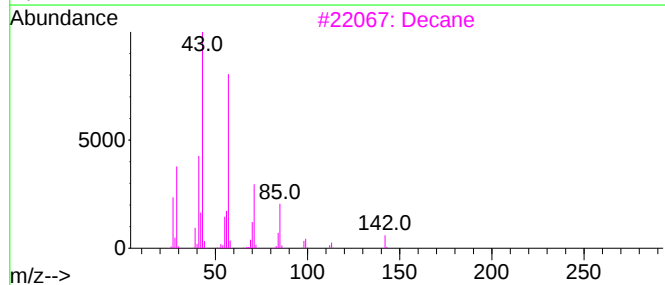
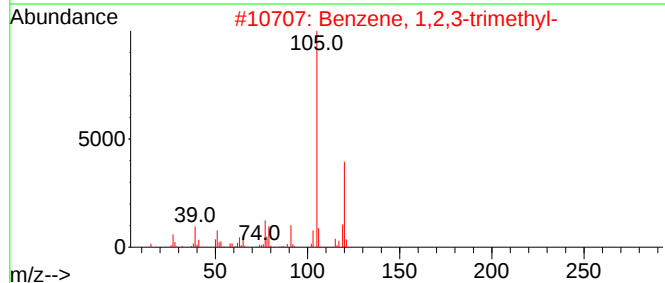
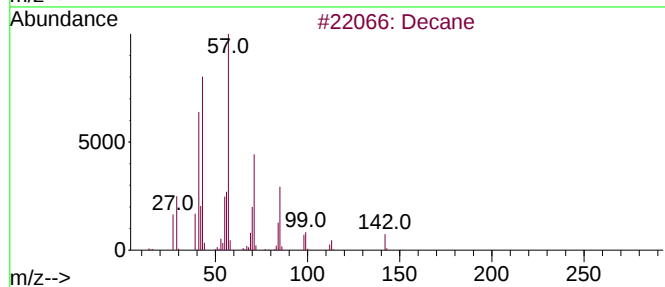
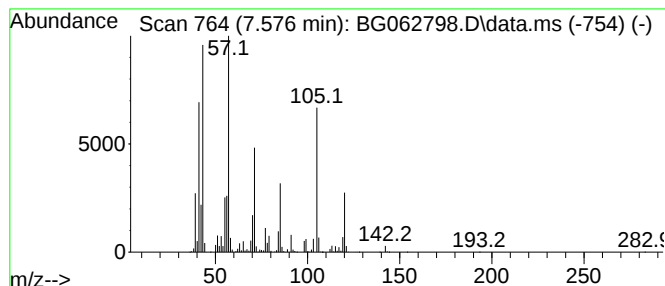
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.576	97.55 ng/ul	36978600	1,4-Dichlorobenzene-d4	7.917

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane	142	C10H22	000124-18-5	93
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	59
3		Decane	142	C10H22	000124-18-5	58
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	56
5		Mesitylene	120	C9H12	000108-67-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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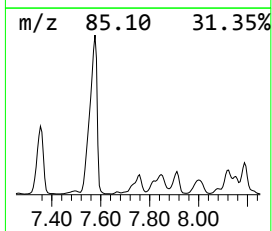
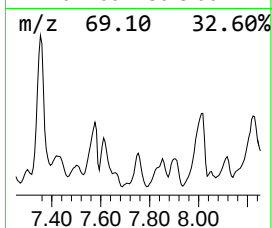
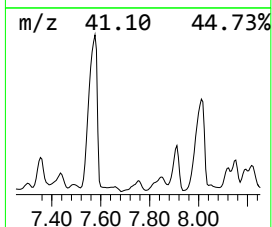
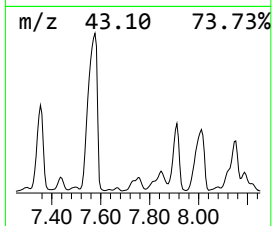
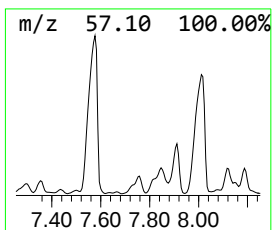
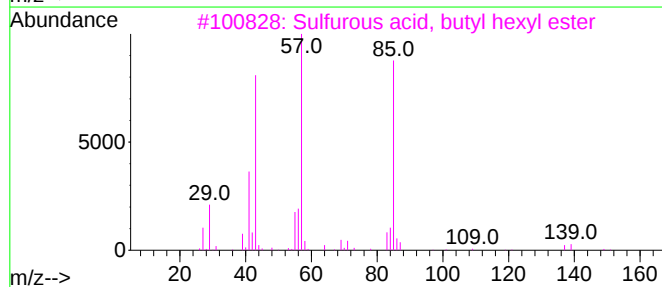
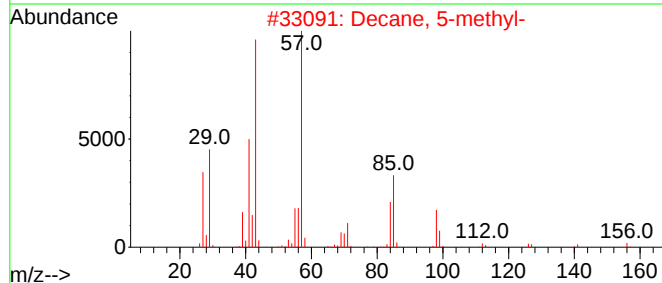
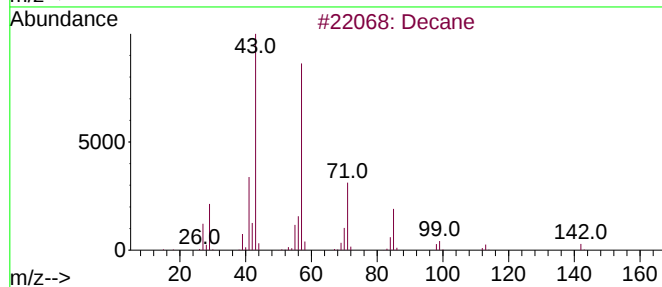
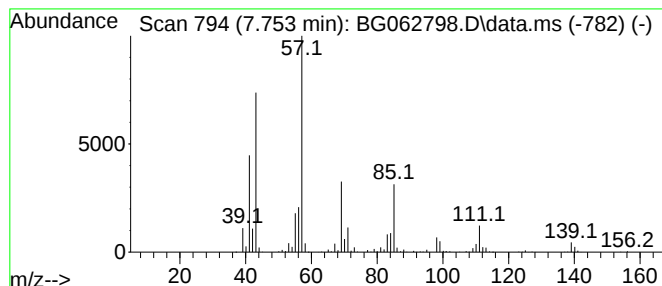
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.753	8.04 ng/ul	3047020	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	53
2			Decane, 5-methyl-	156	C11H24	013151-35-4	53
3			Sulfurous acid, butyl hexyl ester	222	C10H22O3S	1000309-17-3	50
4			Ether, heptyl hexyl	200	C13H28O	007289-40-9	47
5			Hexane, 3-ethyl-3-methyl-	128	C9H20	003074-76-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
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Instrument :
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ClientSampleId :
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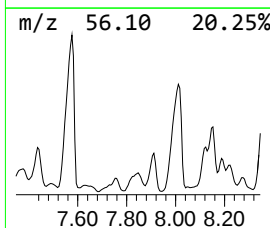
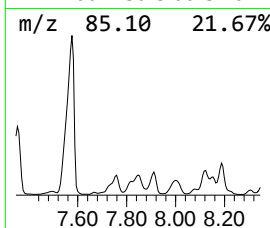
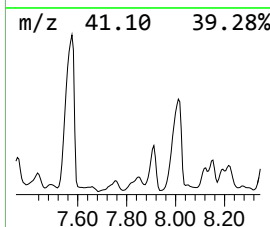
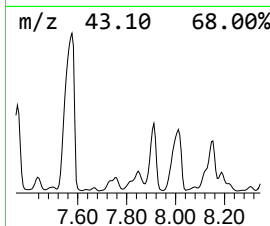
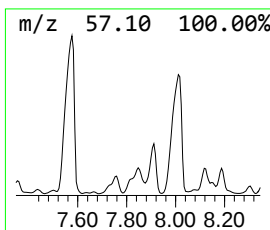
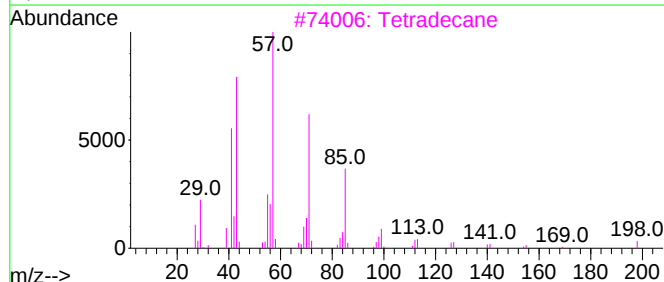
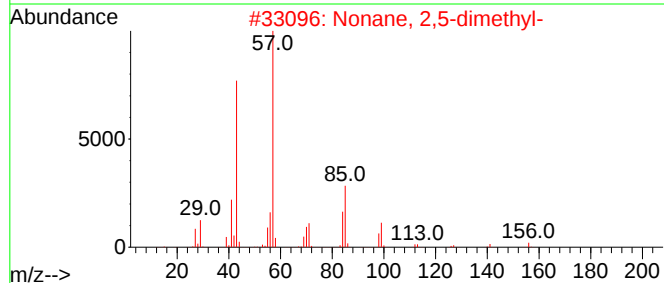
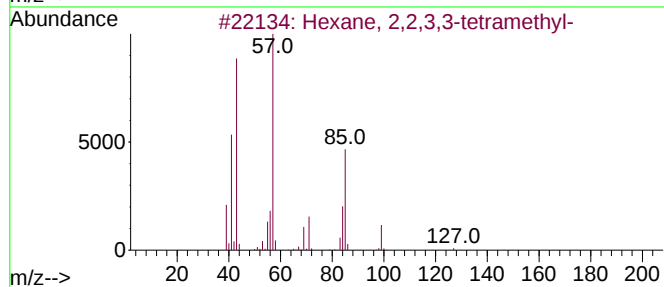
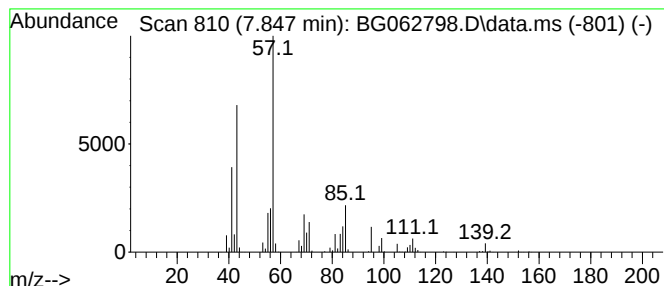
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.847	8.89 ng/ul	3368250	1,4-Dichlorobenzene-d4	7.917

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	53
2		Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	53
3		Tetradecane	198	C14H30	000629-59-4	53
4		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	50
5		Oxalic acid, isobutyl hexyl ester	230	C12H22O4	1000309-37-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
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Instrument :
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ClientSampleId :
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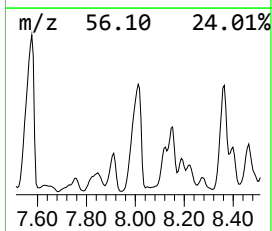
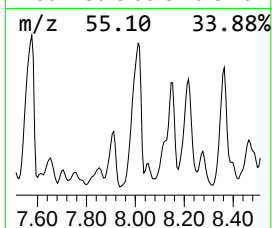
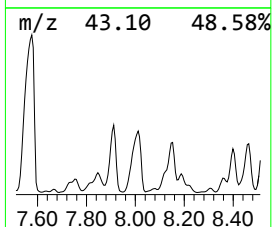
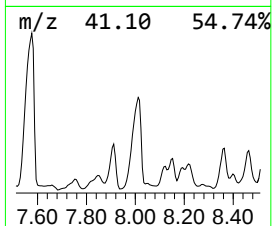
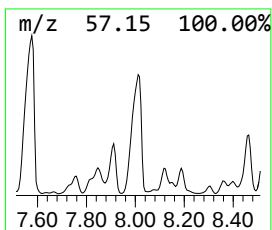
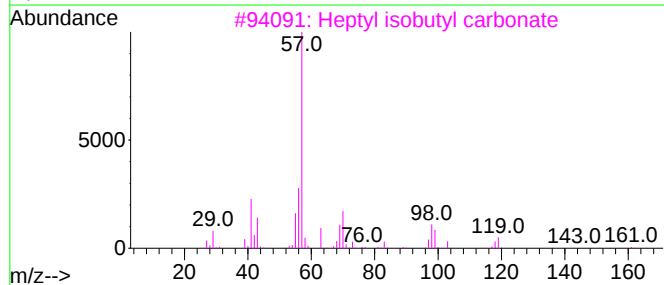
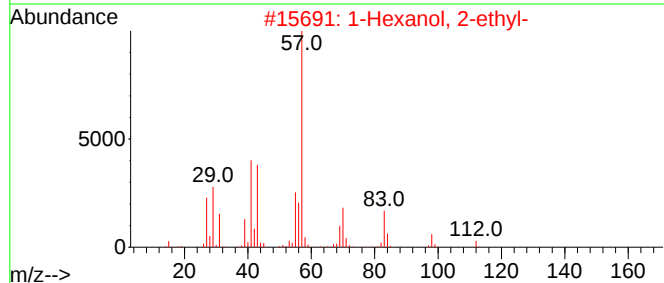
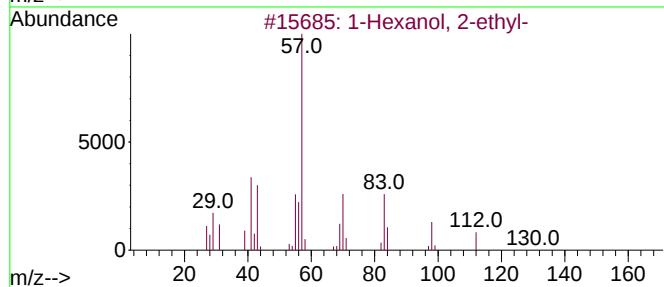
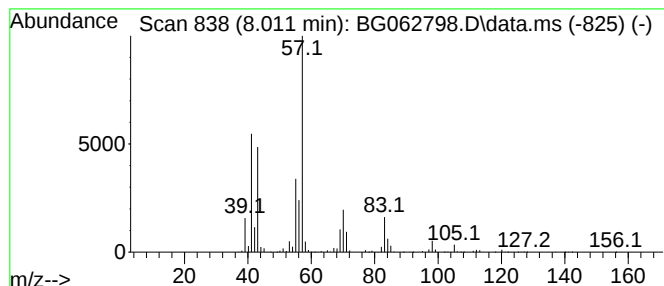
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.011	52.99 ng/ul	20087600	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
3			Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	59
4			Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	53
5			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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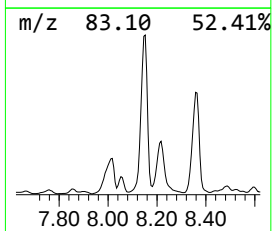
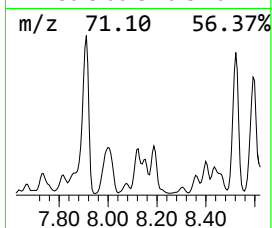
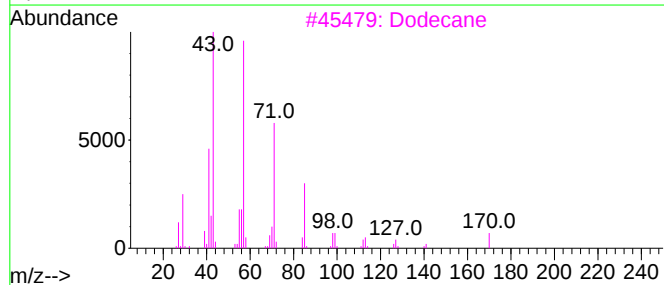
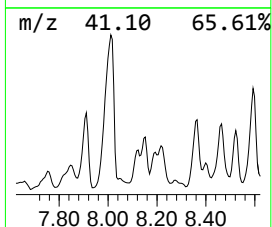
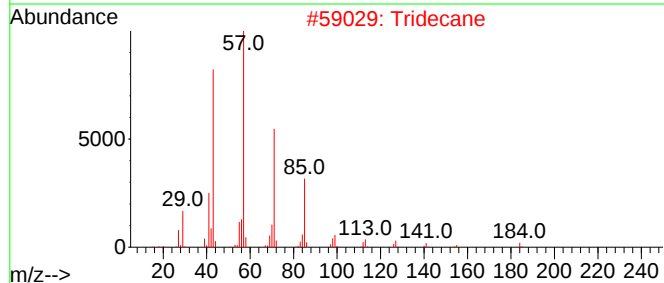
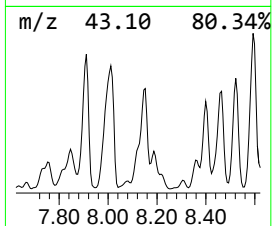
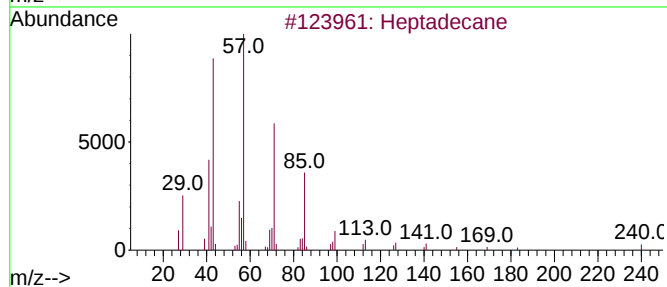
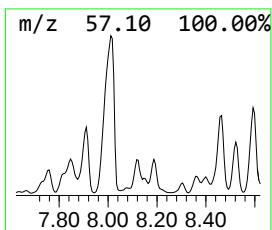
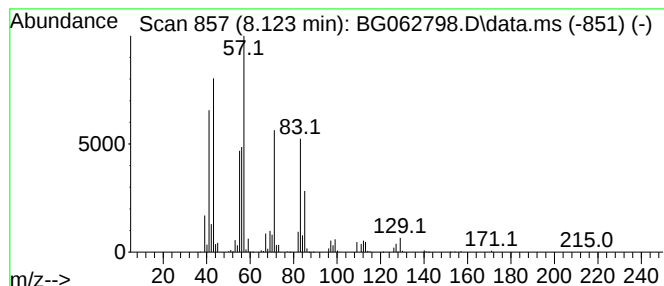
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.123	7.95 ng/ul	3014070	1,4-Dichlorobenzene-d4	7.917

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	46
2		Tridecane	184	C13H28	000629-50-5	46
3		Dodecane	170	C12H26	000112-40-3	46
4		Hexadecane	226	C16H34	000544-76-3	46
5		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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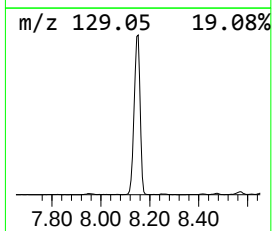
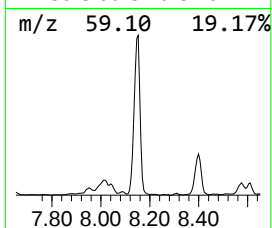
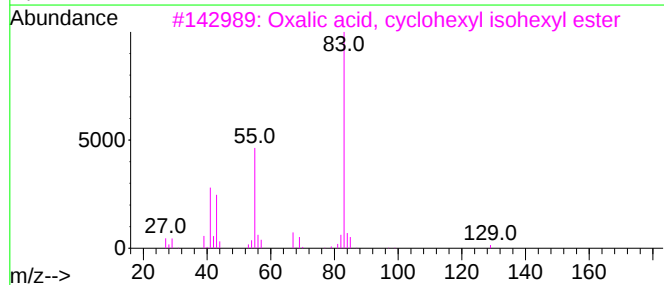
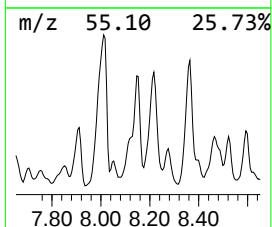
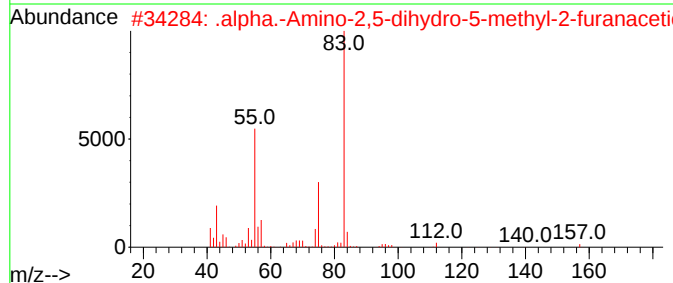
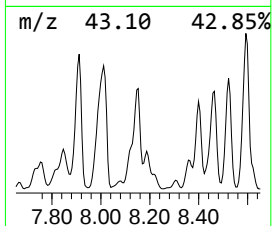
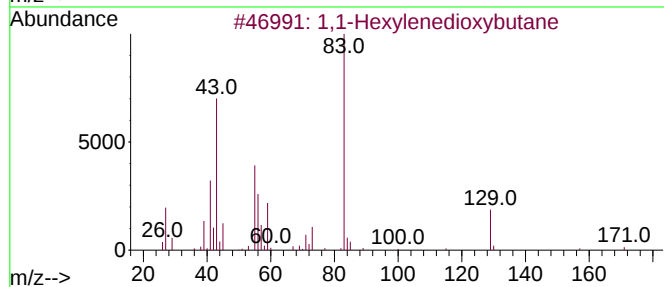
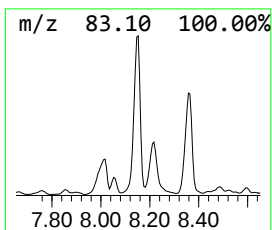
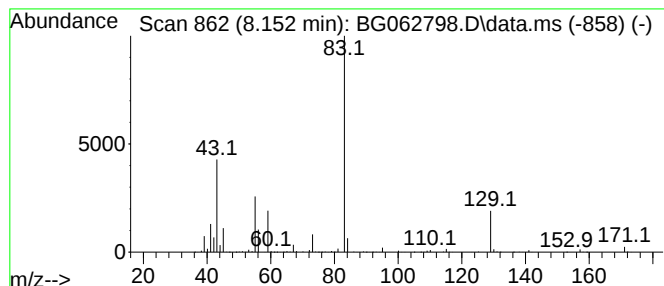
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 1,1-Hexylenedioxybutane Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.152	14.60 ng/ul	5533710	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	56
2			.alpha.-Amino-2,5-dihydro-5-meth...	157	C7H11NO3	018455-25-9	50
3			Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1010309-30-7	43
4			1,3-Cyclohexanedione, 2,2,5,5-te...	168	C10H16O2	000702-50-1	28
5			2-Dimethylamino-4-methyl-pent-4-...	138	C8H14N2	094492-10-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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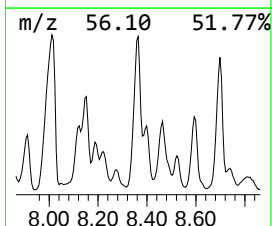
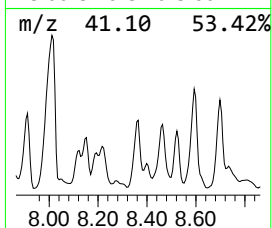
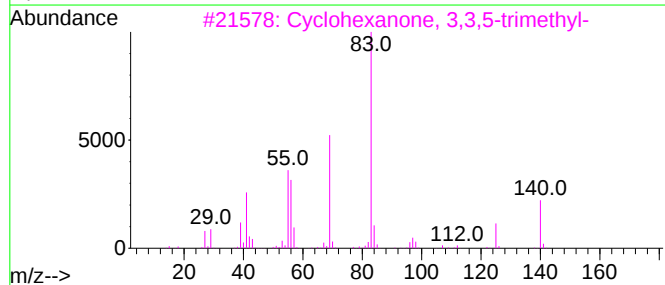
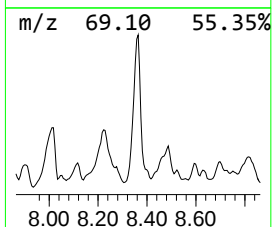
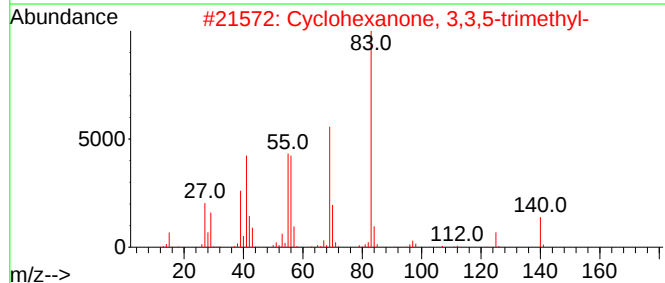
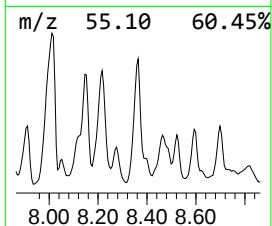
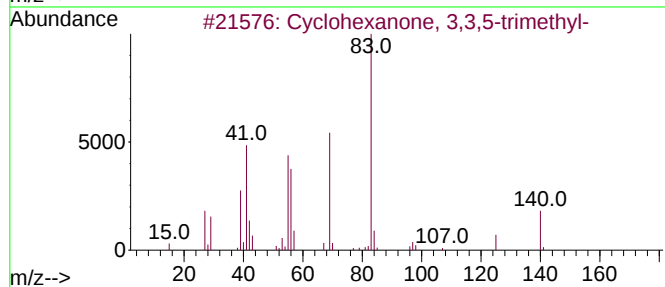
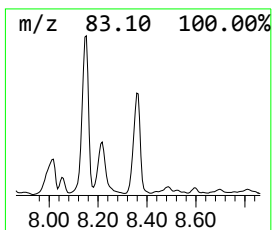
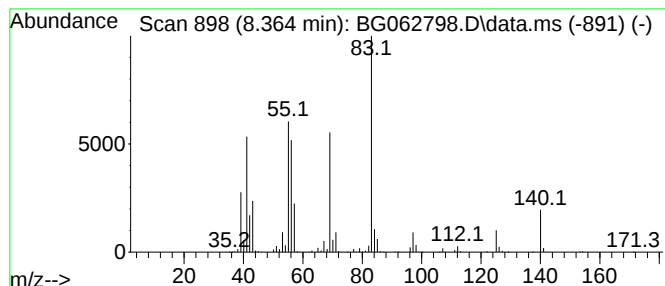
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Cyclohexanone, 3,3,5-trimet... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.364	20.77 ng/ul	7873070	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	81
5			cis-3-Decene	140	C10H20	019398-86-8	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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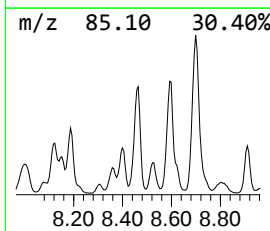
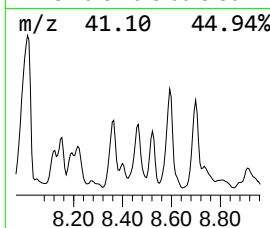
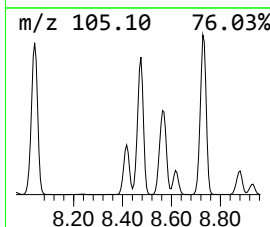
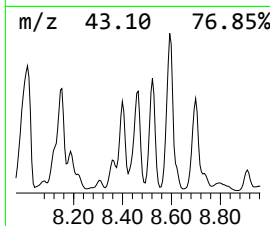
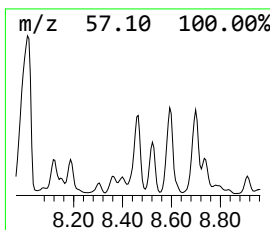
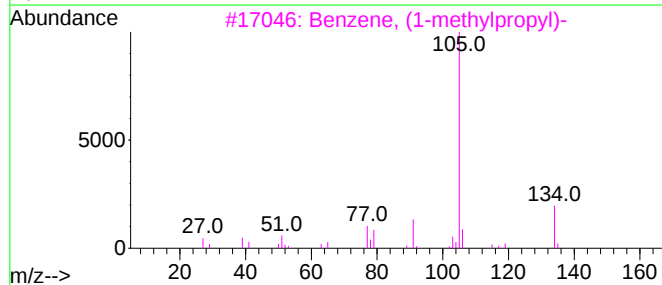
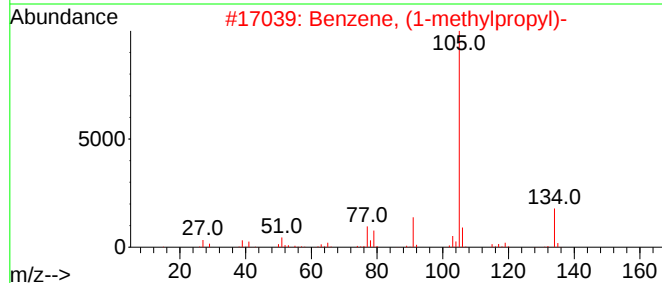
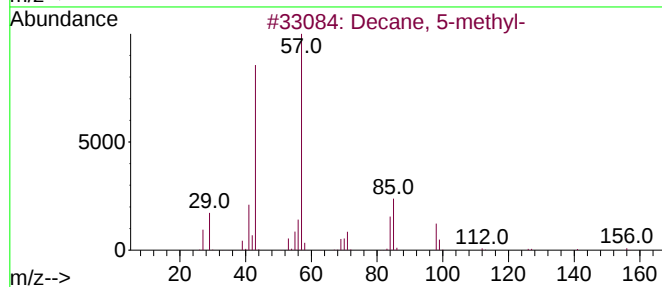
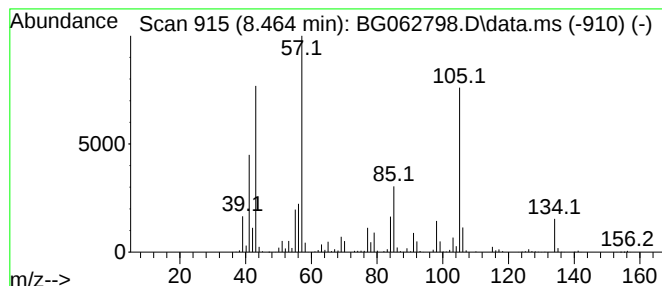
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.464	21.27 ng/ul	8061470	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	46
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38
4			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	38
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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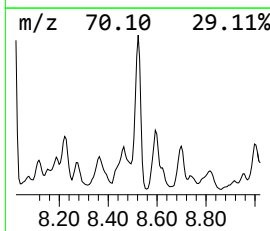
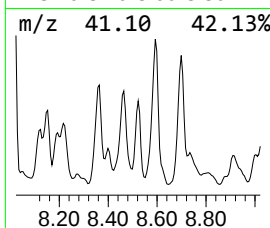
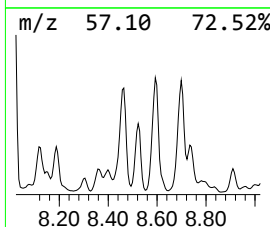
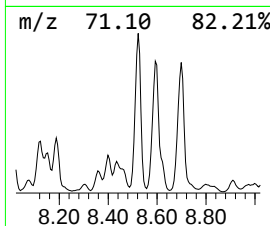
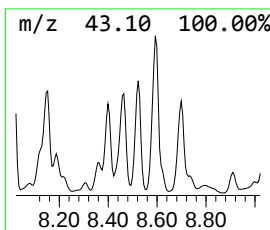
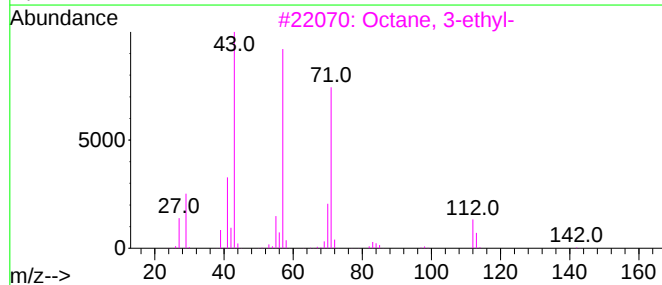
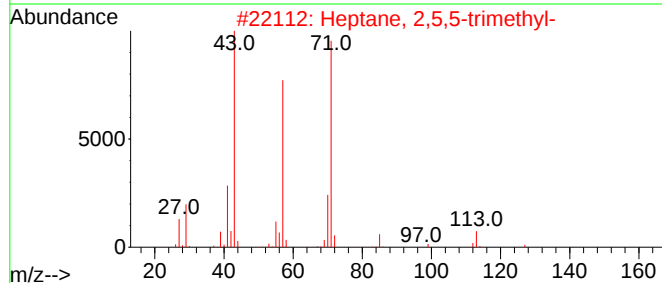
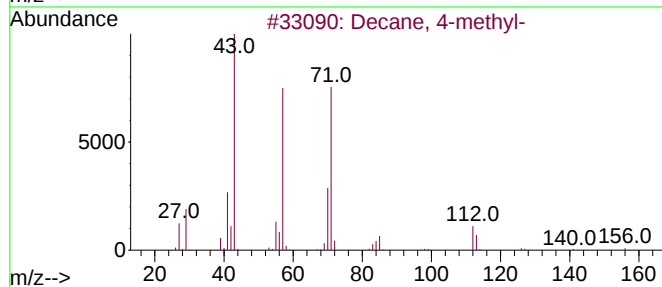
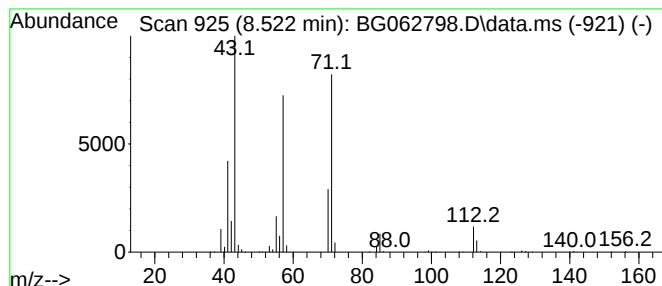
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.522	9.39 ng/ul	3561060	1,4-Dichlorobenzene-d4	7.917

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 4-methyl-	156	C11H24	002847-72-5	91
2		Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	78
3		Octane, 3-ethyl-	142	C10H22	005881-17-4	78
4		Dodecane, 4-methyl-	184	C13H28	006117-97-1	72
5		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
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ClientSampleId :
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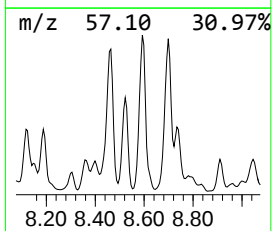
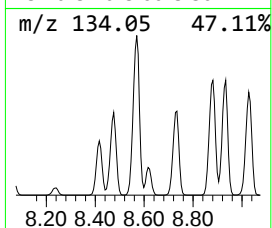
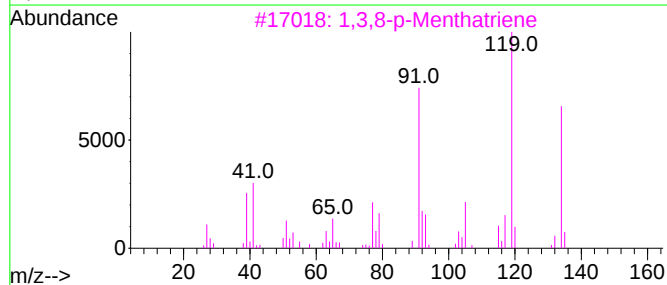
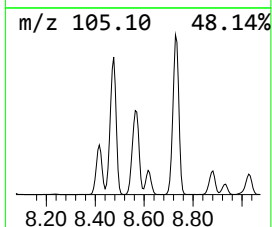
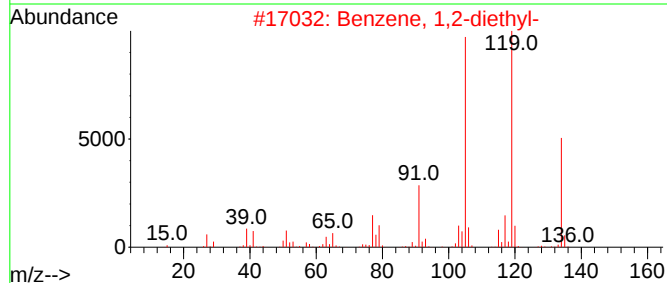
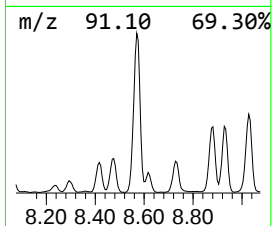
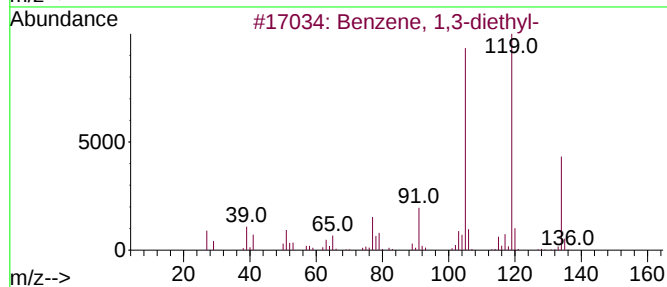
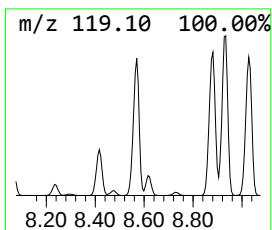
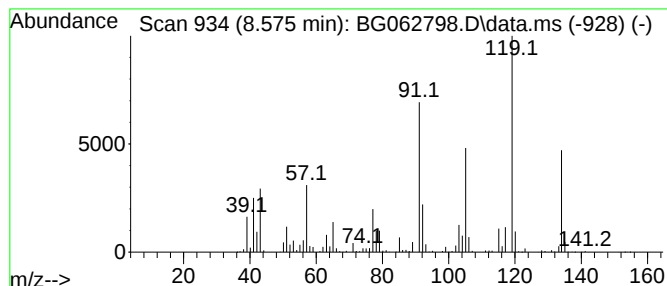
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1,3-diethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.575	17.31 ng/ul	6561300	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	87
3			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	83
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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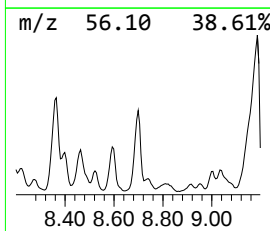
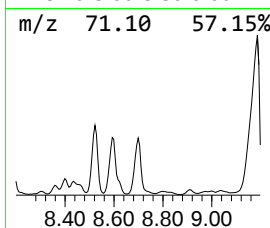
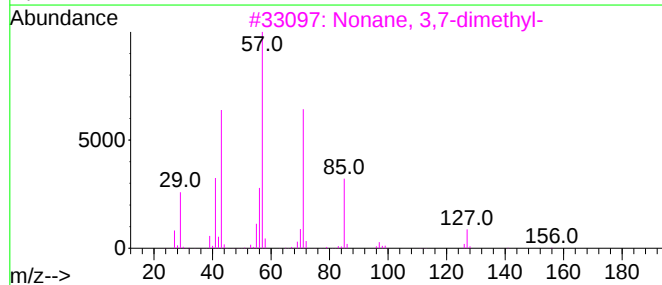
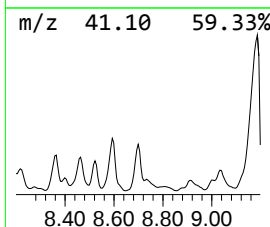
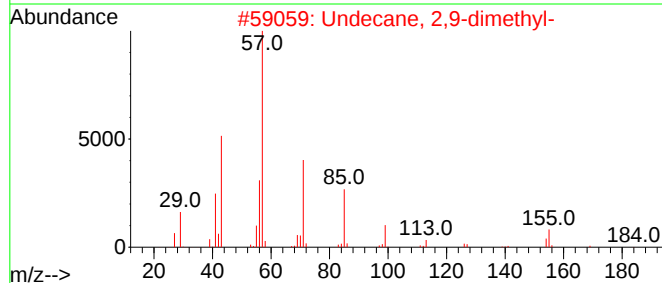
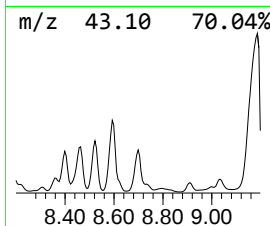
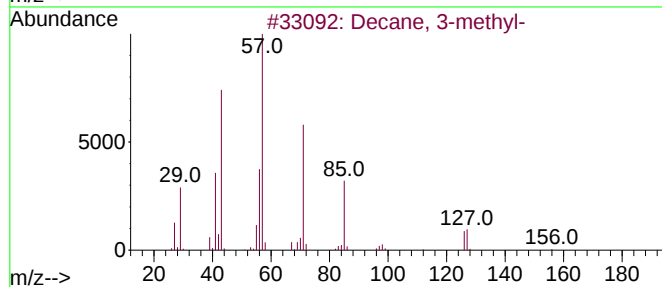
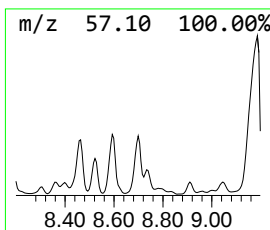
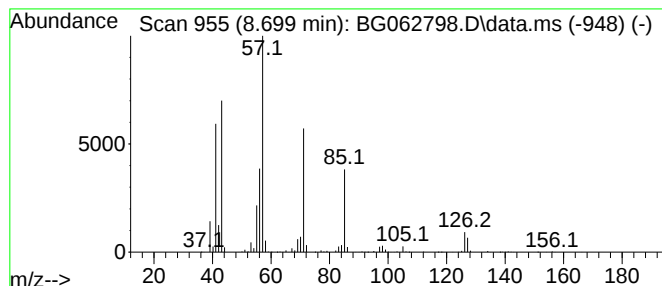
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.699	21.00 ng/ul	7959450	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	90
2			Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	72
3			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	58
4			Decane, 3-methyl-	156	C11H24	013151-34-3	53
5			Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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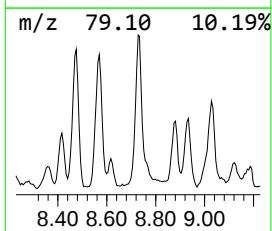
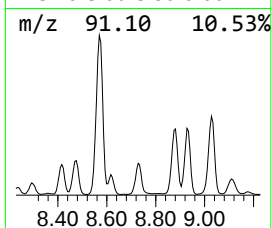
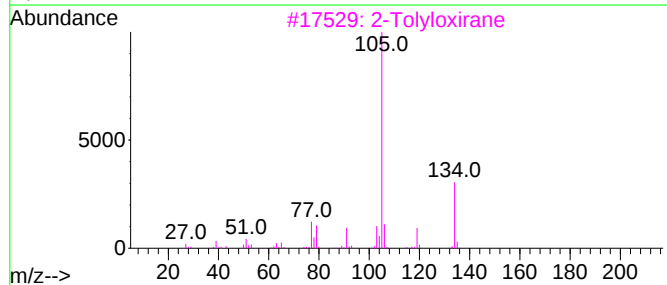
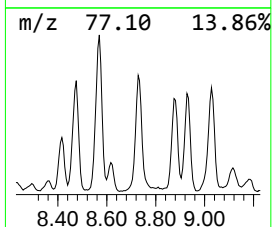
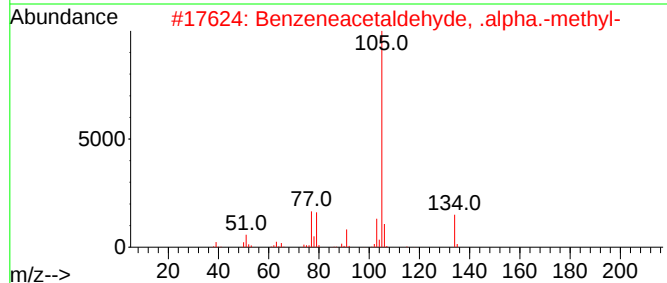
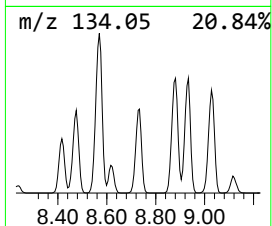
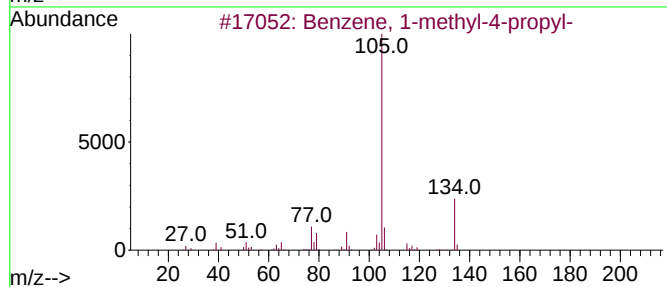
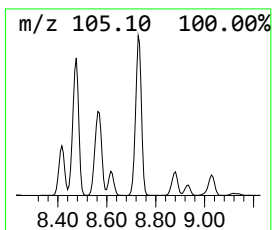
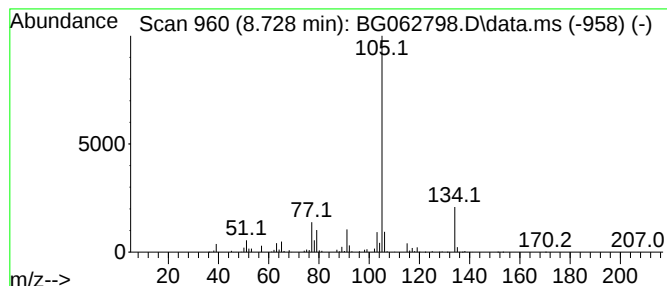
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 Benzene, 1-methyl-4-propyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.728	13.41 ng/ul	5081490	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91
3			2-Tolyloxirane	134	C9H10O	002783-26-8	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
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ClientSampleId :
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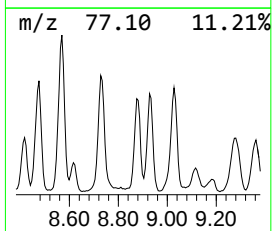
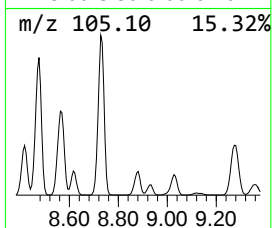
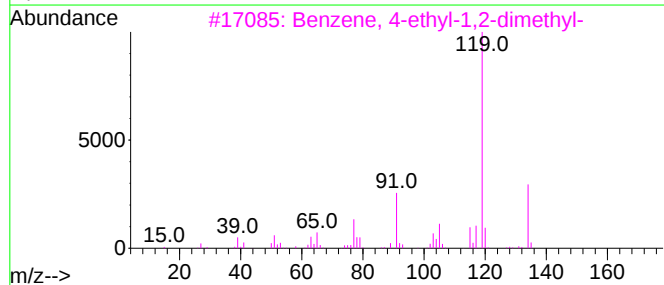
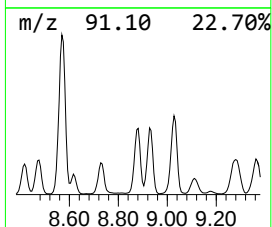
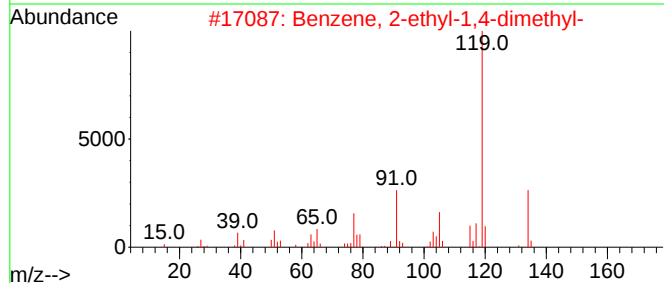
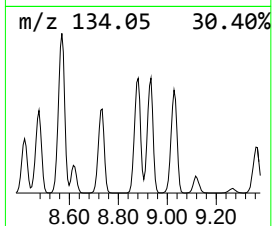
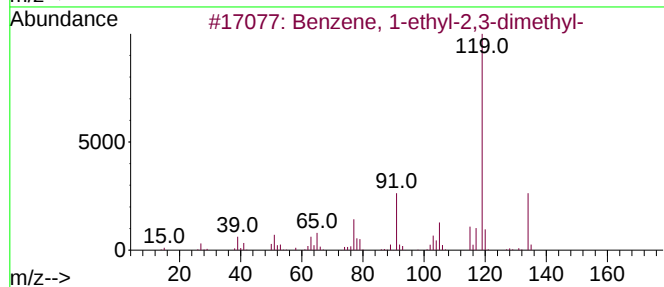
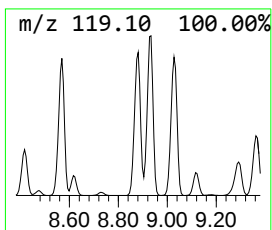
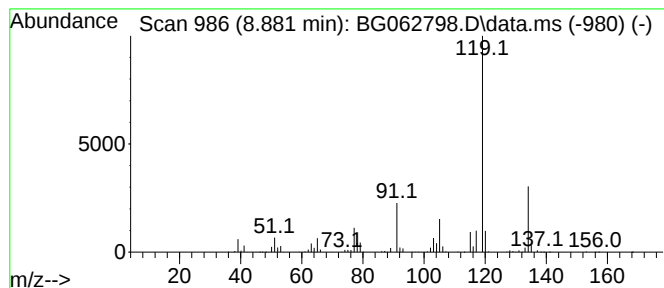
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.881	11.85 ng/ul	4490620	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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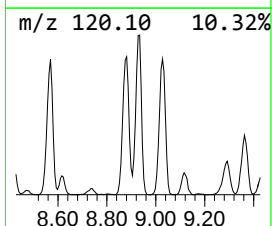
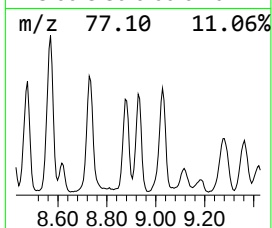
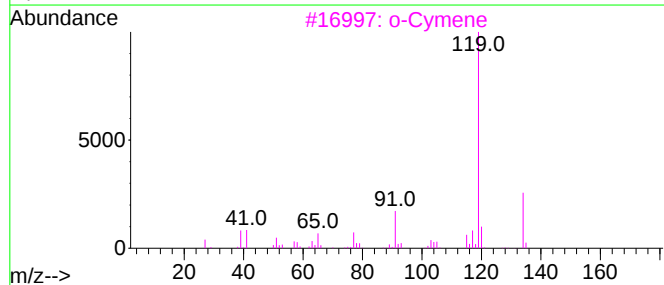
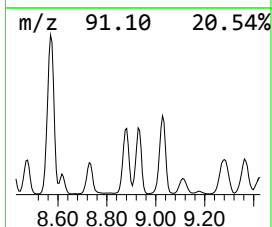
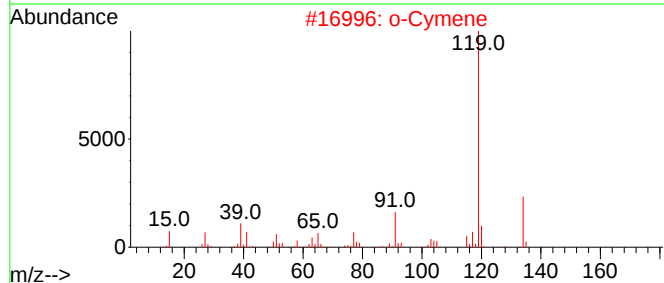
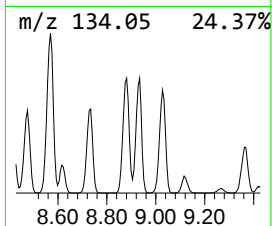
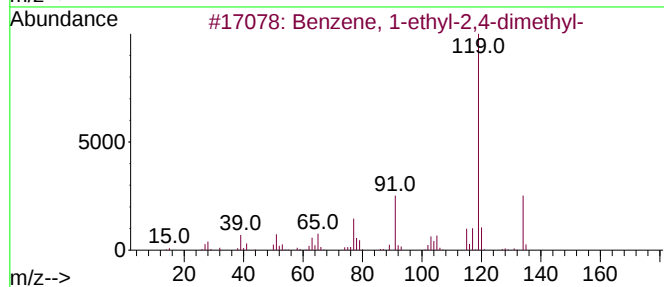
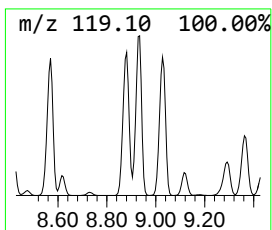
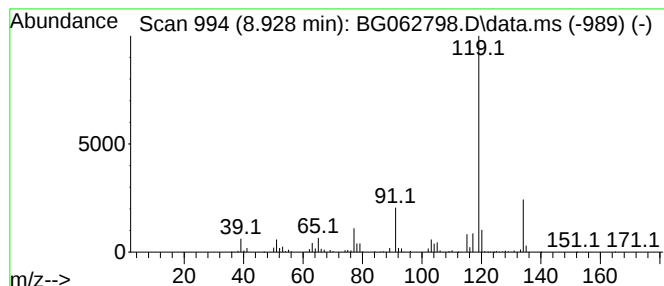
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	17.73 ng/ul	6721870	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			o-Cymene	134	C10H14	000527-84-4	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

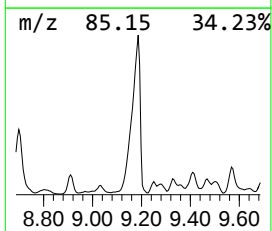
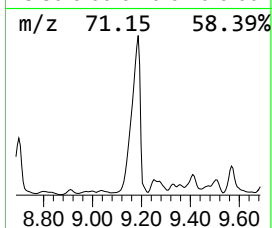
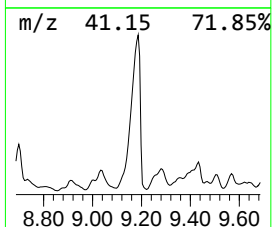
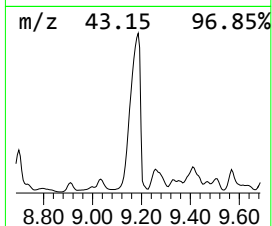
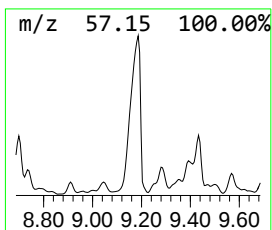
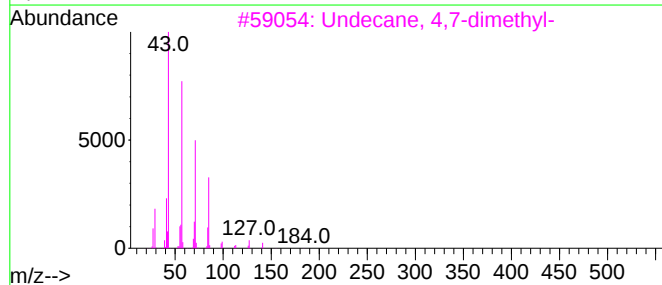
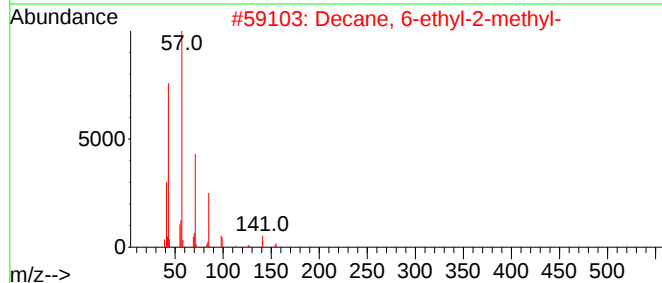
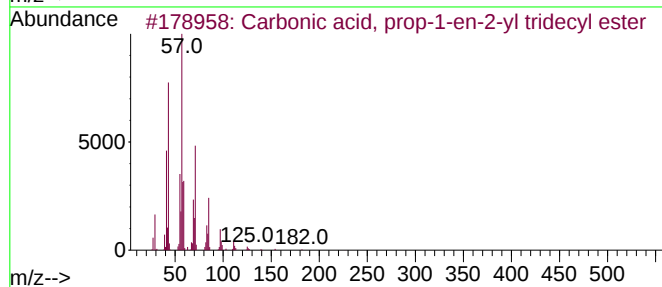
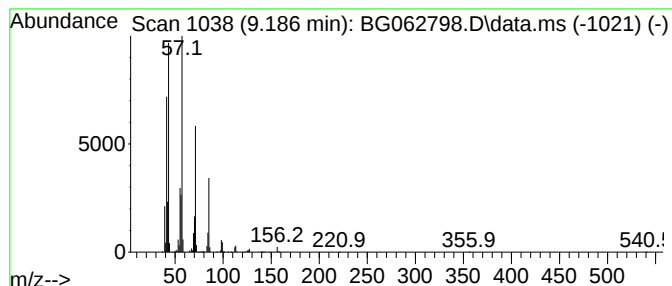
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Carbonic acid, prop-1-en-2-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.186	101.64 ng/ul	38528200	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	86
2			Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	59
3			Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	59
4			2-Bromononane	206	C9H19Br	002216-35-5	53
5			Decane	142	C10H22	000124-18-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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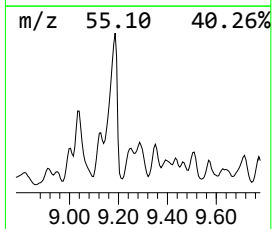
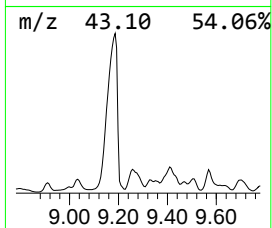
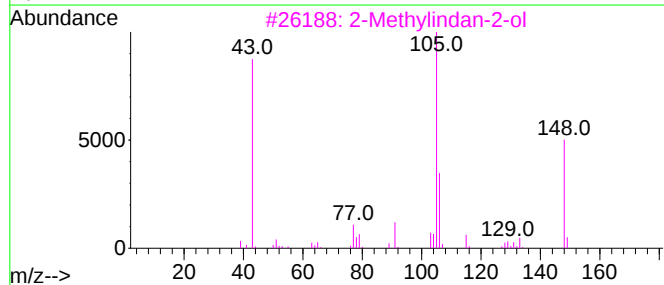
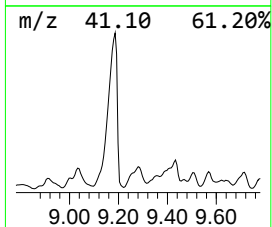
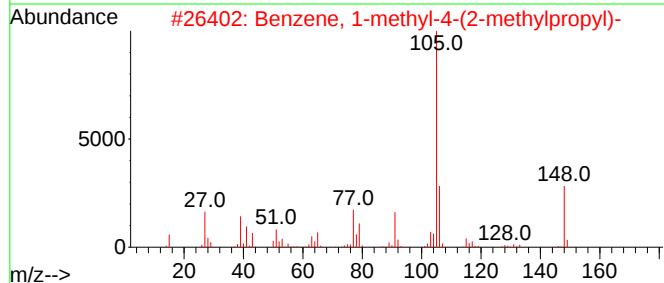
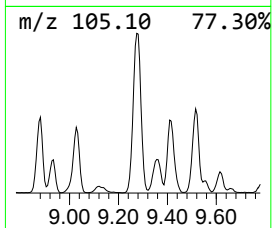
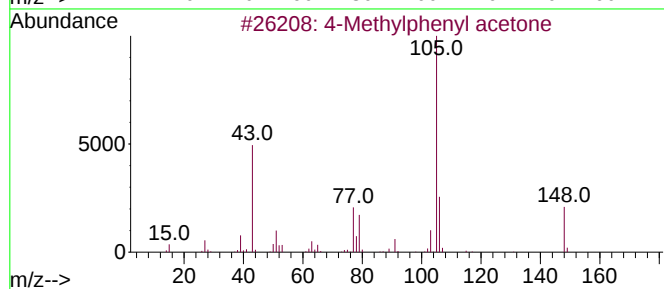
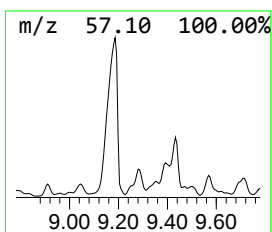
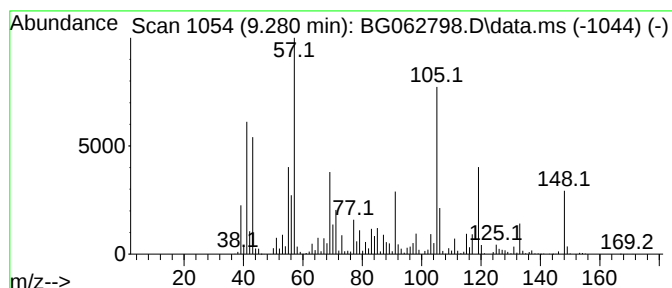
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.280	29.06 ng/ul	11014300	1,4-Dichlorobenzene-d4	7.917

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylphenyl acetone	148	C10H12O	002096-86-8	42
2			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	38
3			2-Methylindan-2-ol	148	C10H12O	033223-84-6	30
4			3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	25
5			Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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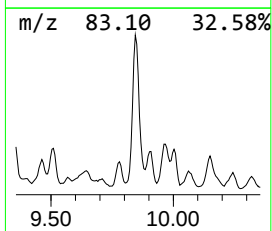
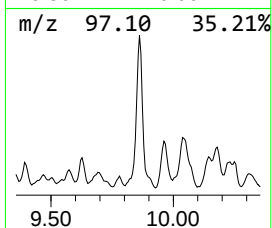
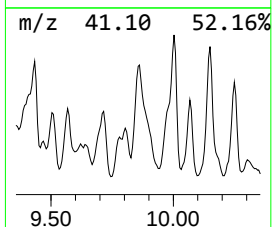
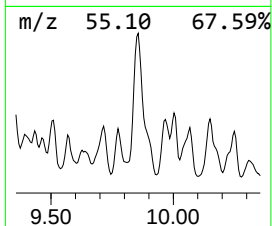
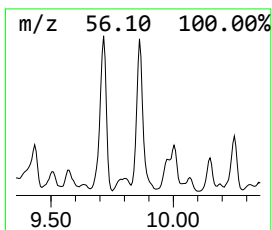
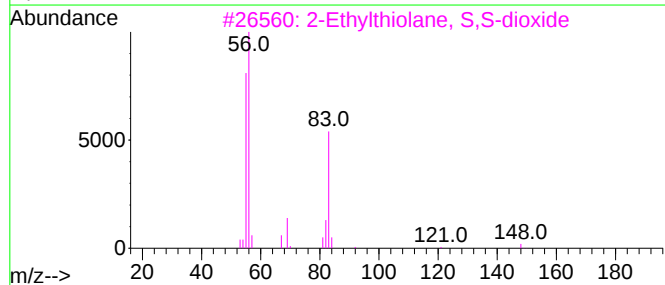
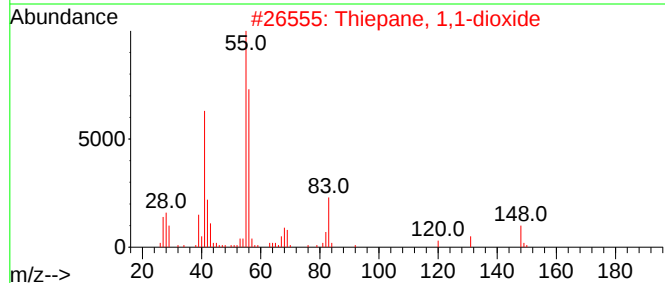
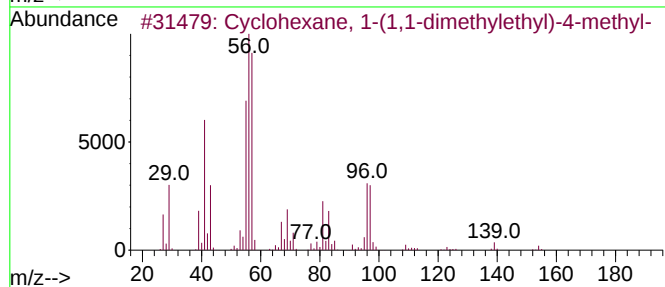
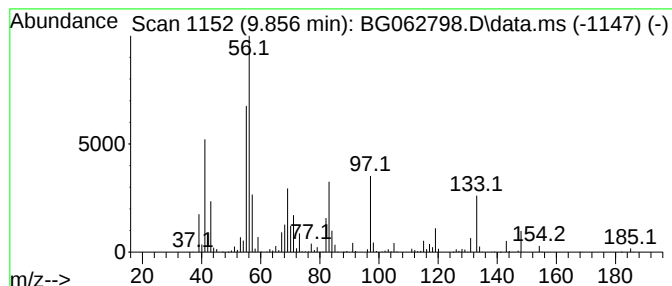
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 (DEL) Alkane: Cyclic9.856 Concentration Rank 68

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.856	3.61 ng/ul	8195450	Naphthalene-d8	10.714

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-(1,1-dimethylethy...	154	C11H22	075736-66-2	47
2			Thiepane, 1,1-dioxide	148	C6H12O2S	006251-33-8	47
3			2-Ethylthiolane, S,S-dioxide	148	C6H12O2S	010178-59-3	43
4			1-Decene, 2-methyl-	154	C11H22	013151-27-4	38
5			Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

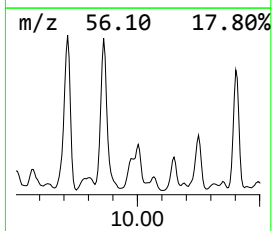
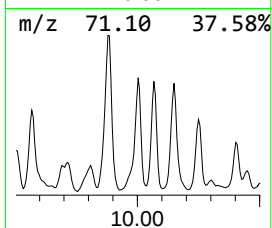
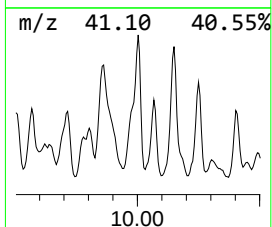
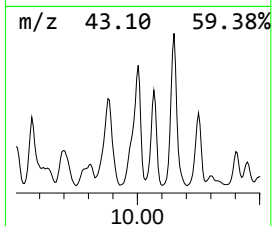
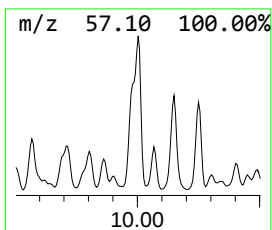
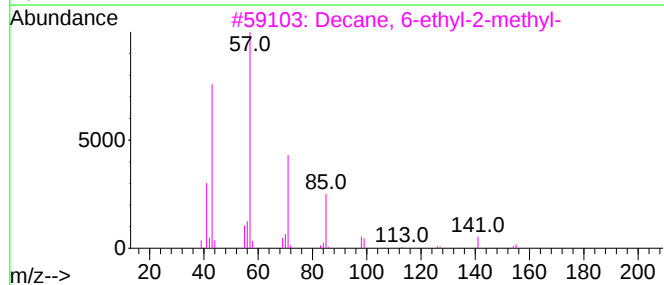
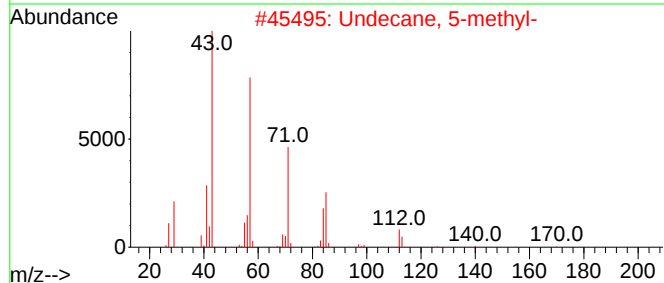
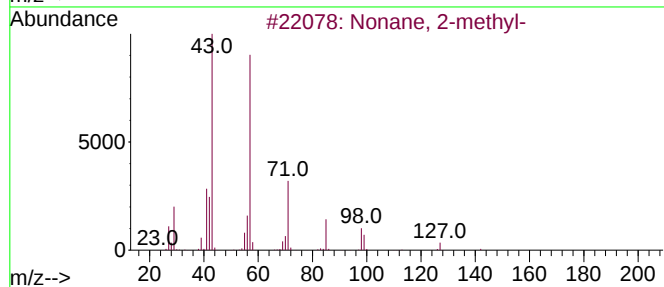
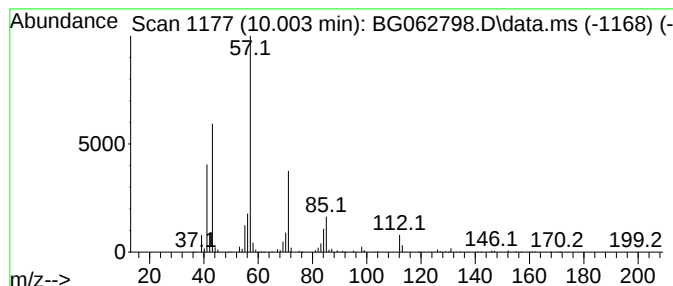
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.003	3.85 ng/ul	8734640	Naphthalene-d8	10.714	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-	142	C10H22	000871-83-0	64
2	Undecane, 5-methyl-	170	C12H26	001632-70-8	64
3	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	59
4	Decane, 2,6,6-trimethyl-	184	C13H28	062108-24-1	53
5	Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

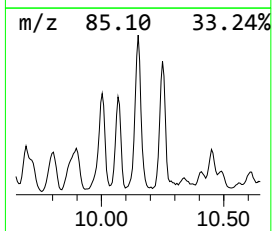
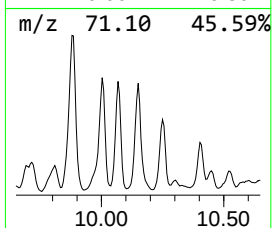
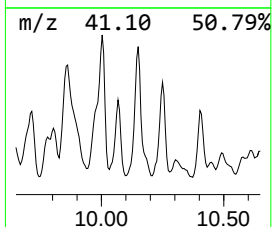
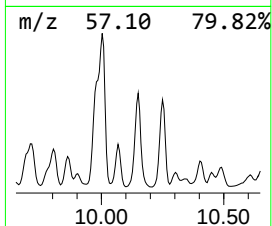
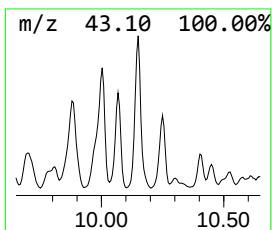
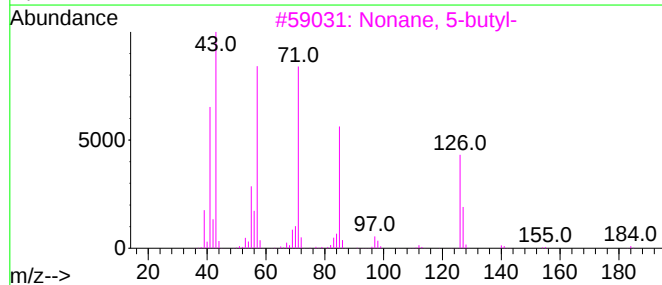
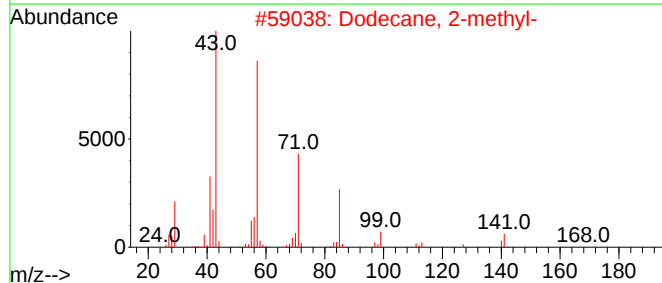
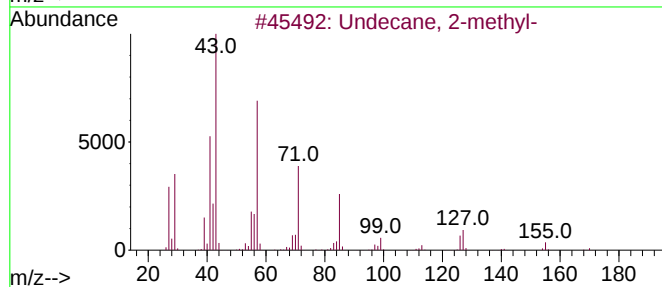
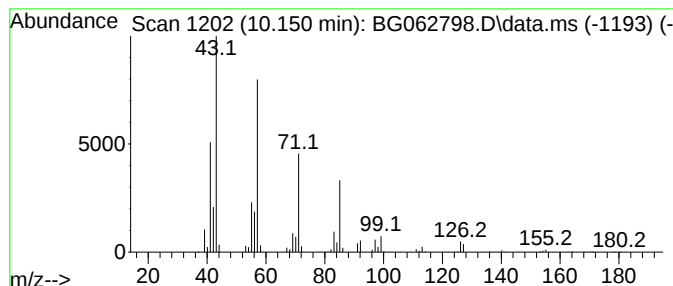
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.150	3.69 ng/ul	8378390	Naphthalene-d8	10.714	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2-methyl-	170	C12H26	007045-71-8	78
2	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
3	Nonane, 5-butyl-	184	C13H28	017312-63-9	64
4	Octacosane	394	C28H58	000630-02-4	64
5	Heptadecane	240	C17H36	000629-78-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

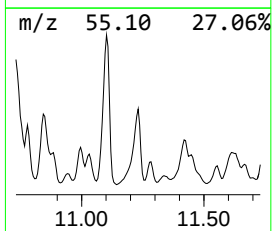
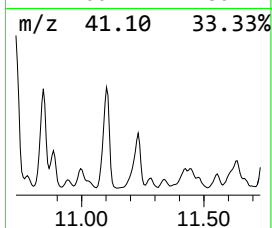
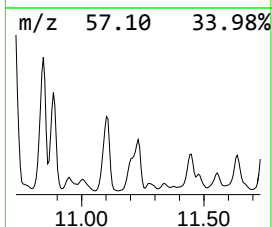
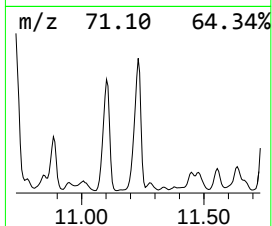
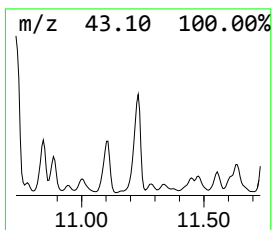
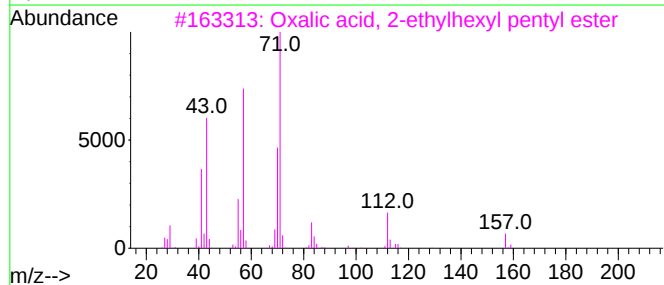
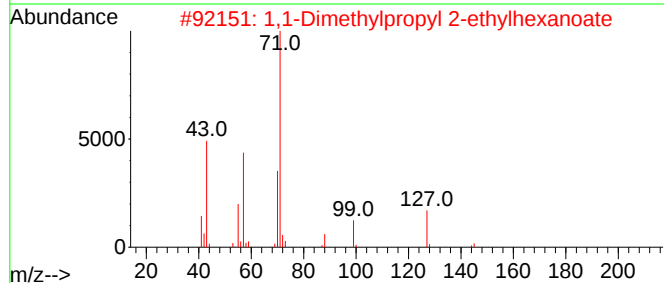
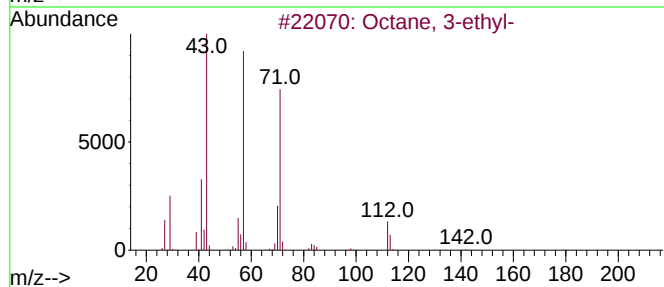
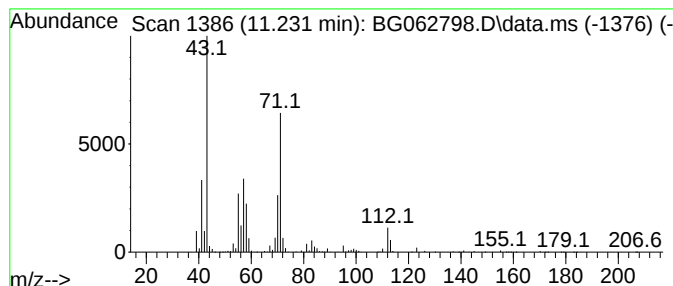
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.231	4.90 ng/ul	11120900	Naphthalene-d8	10.714	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-	142	C10H22	005881-17-4	59
2	1,1-Dimethylpropyl 2-ethylhexanoate	214	C13H26O2	1000099-99-2	50
3	Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	47
4	Sulfurous acid, nonyl 2-pentyl e...	278	C14H30O3S	1010309-15-8	47
5	Decane, 4-methyl-	156	C11H24	002847-72-5	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

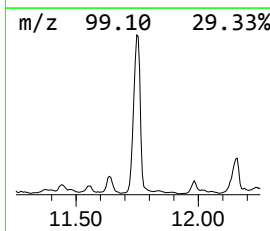
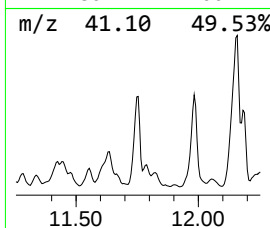
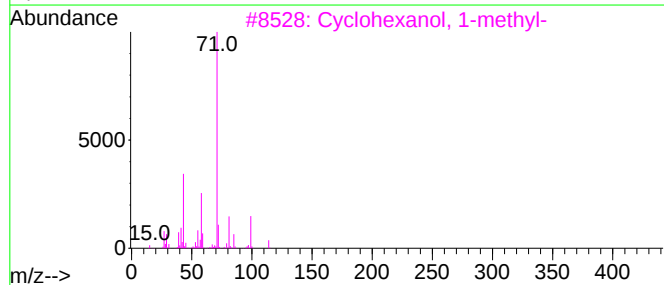
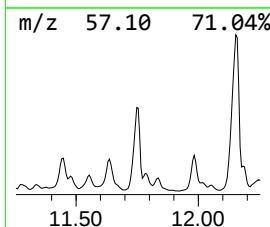
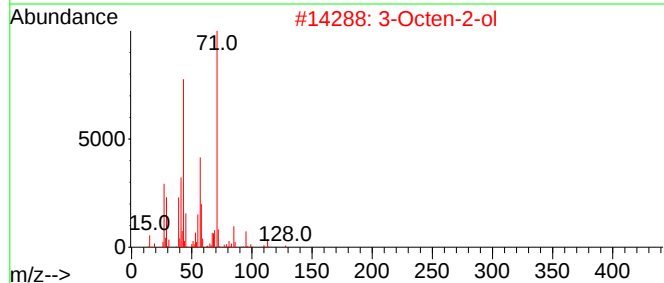
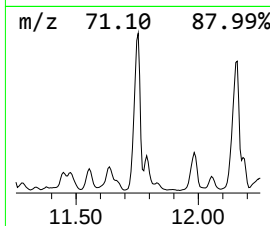
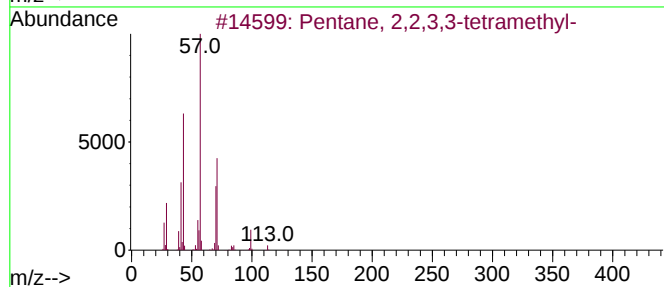
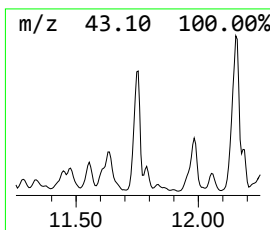
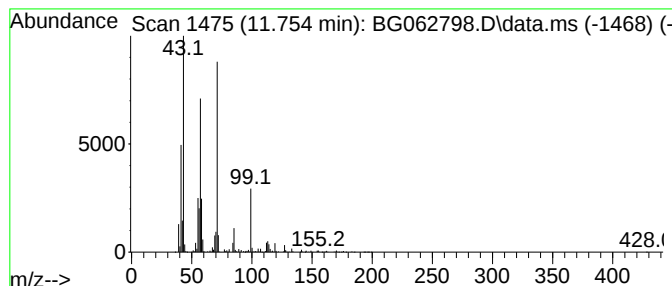
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.754	5.11 ng/ul	11605500	Naphthalene-d8	10.714	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	64
2	3-Octen-2-ol	128	C8H16O	076649-14-4	52
3	Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	52
4	2,4-Pentanedione, 3-butyl-	156	C9H16O2	001540-36-9	50
5	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

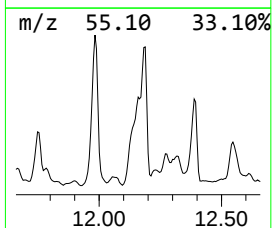
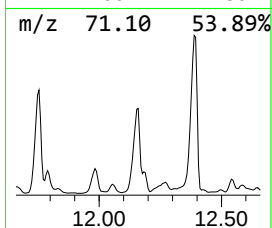
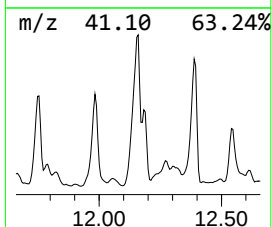
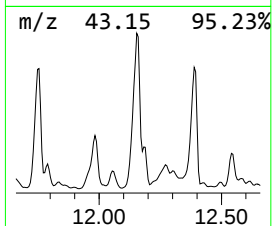
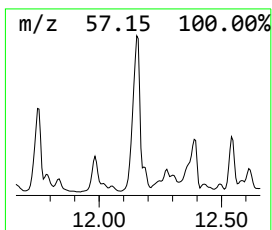
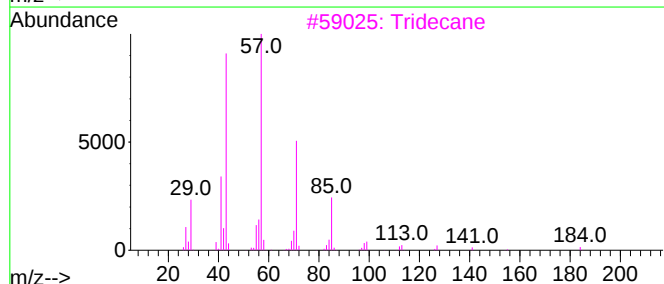
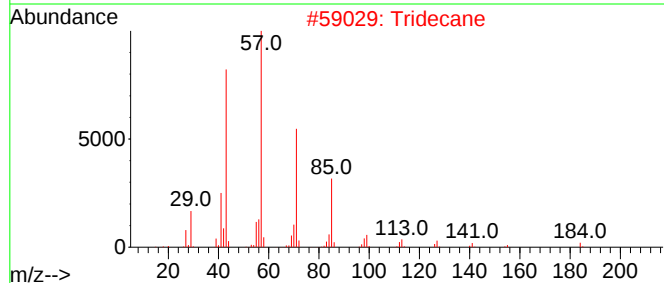
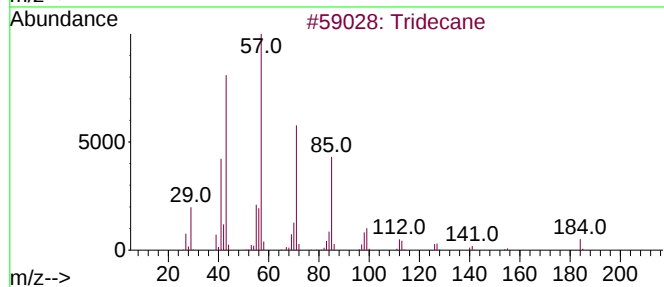
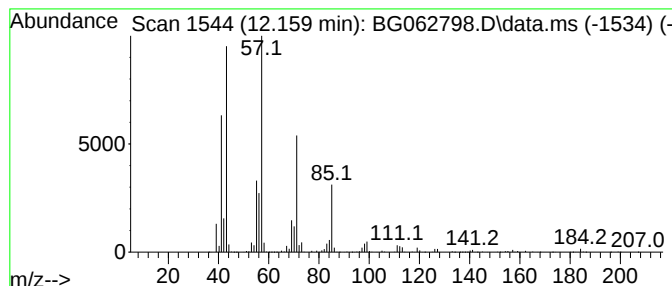
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.159	7.87 ng/ul	17867700	Naphthalene-d8	10.714

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	93
2		Tridecane	184	C13H28	000629-50-5	87
3		Tridecane	184	C13H28	000629-50-5	81
4		Tetradecane	198	C14H30	000629-59-4	80
5		Dodecane	170	C12H26	000112-40-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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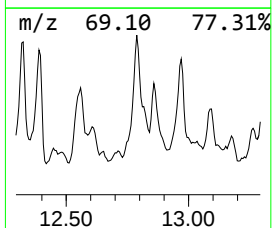
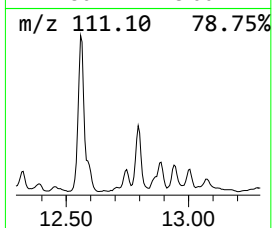
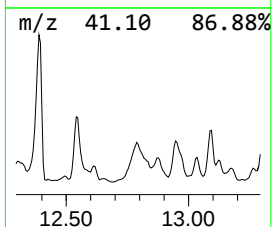
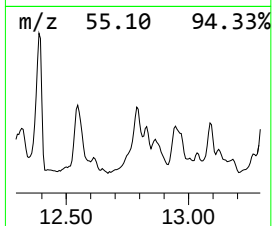
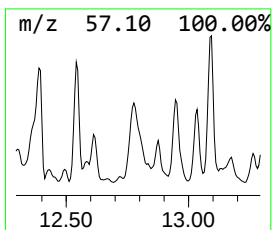
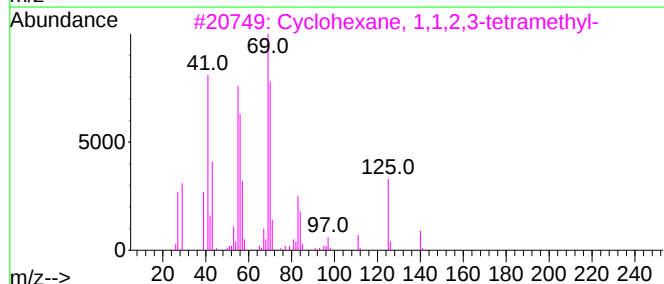
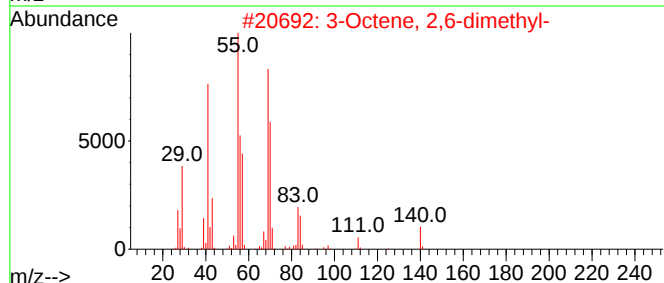
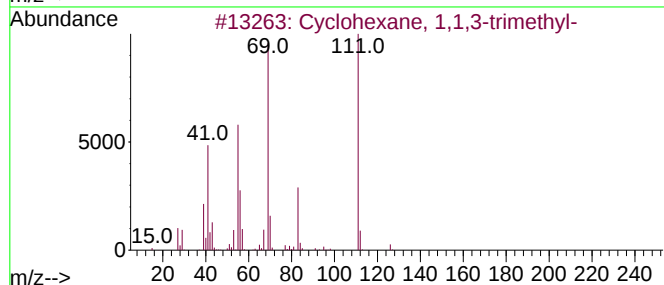
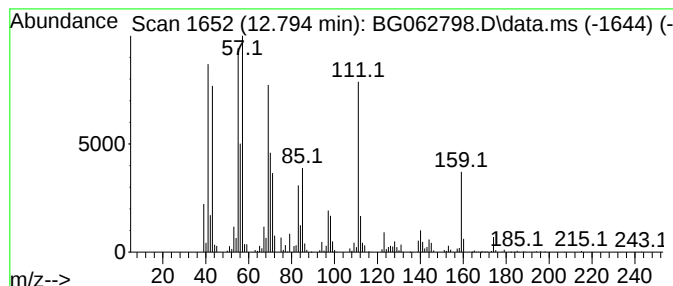
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 (DEL) Alkane: Cyclic12.794 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.794	16.18 ng/ul	4768820	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	46
2			3-Octene, 2,6-dimethyl-	140	C10H20	006874-28-8	45
3			Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	45
4			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	43
5			Cyclopentane, ethyl-	98	C7H14	001640-89-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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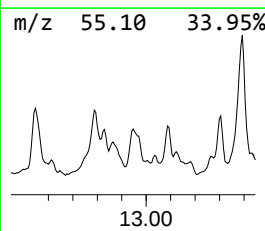
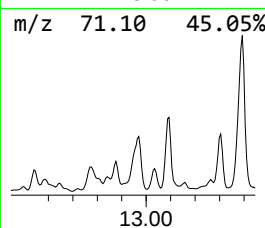
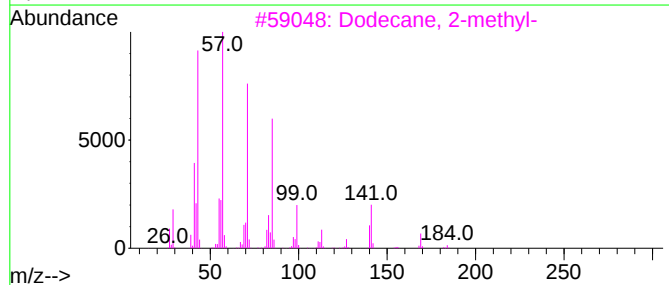
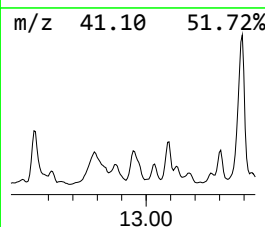
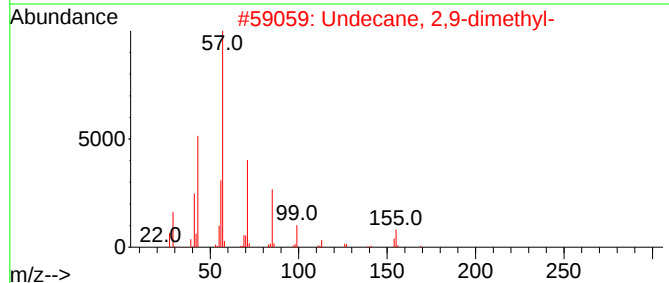
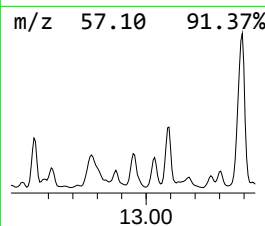
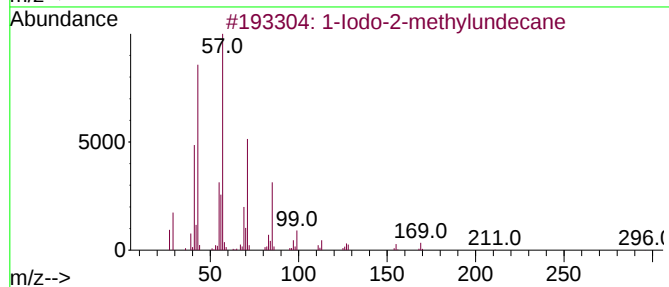
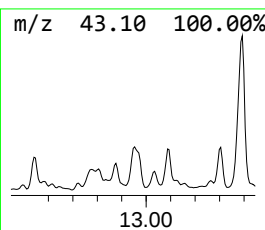
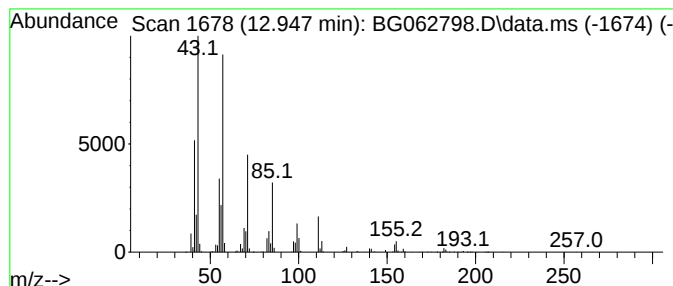
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 1-Iodo-2-methylundecane Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.947	15.35 ng/ul	4523300	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	80
2			Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	78
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
4			Tridecane, 2-methyl-	198	C14H30	001560-96-9	72
5			Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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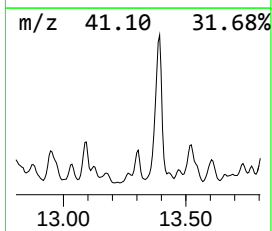
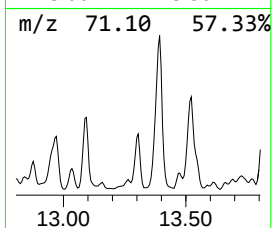
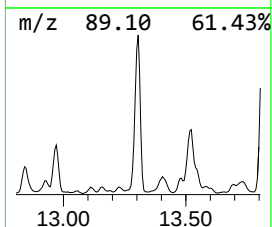
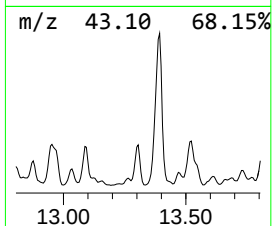
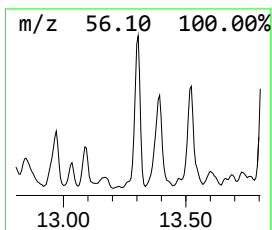
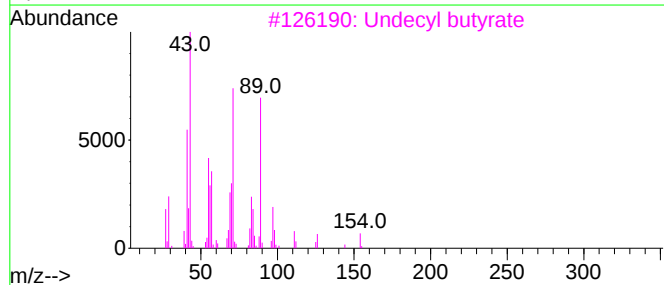
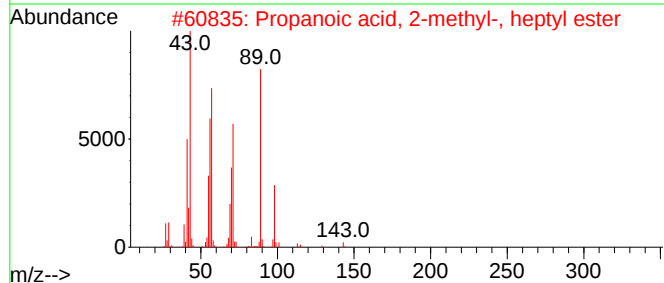
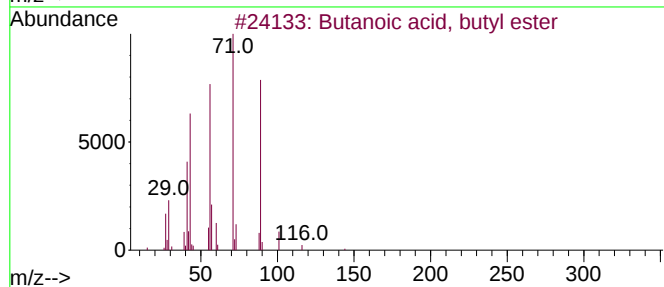
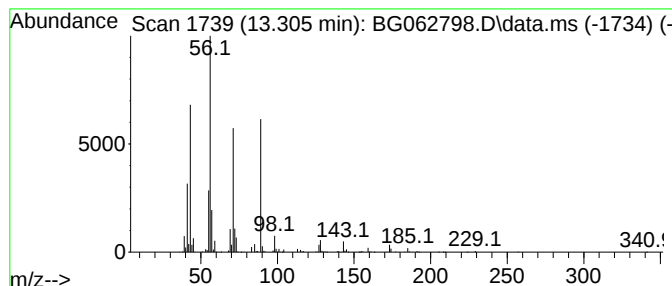
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 unknown-02 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.305	13.06 ng/ul	3847550	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	47
2			Propanoic acid, 2-methyl-, heptyl...	186	C11H22O2	002349-13-5	38
3			Undecyl butyrate	242	C15H30O2	1000428-74-4	38
4			Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	37
5			Propanoic acid, 2-methyl-, hexyl...	172	C10H20O2	002349-07-7	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

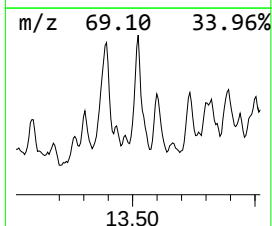
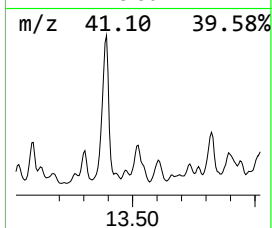
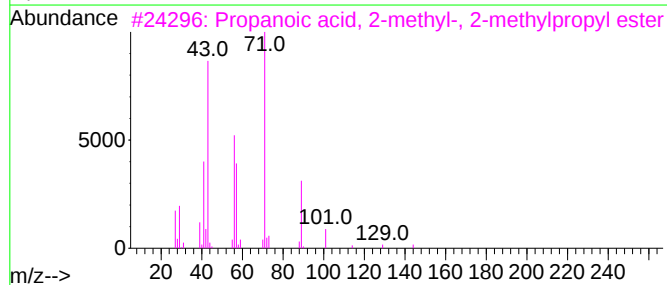
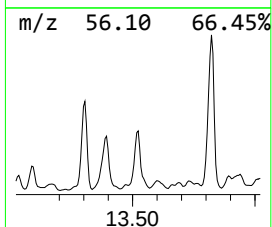
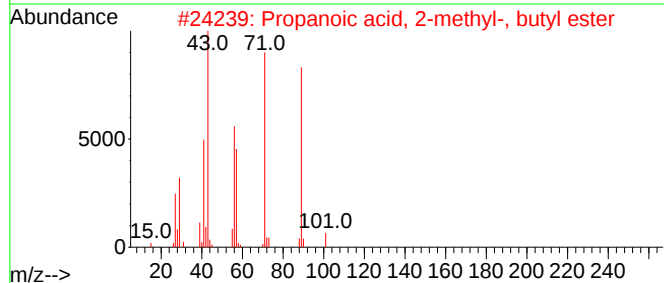
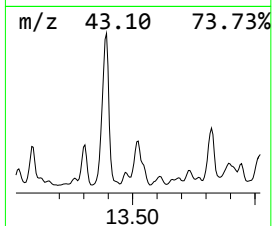
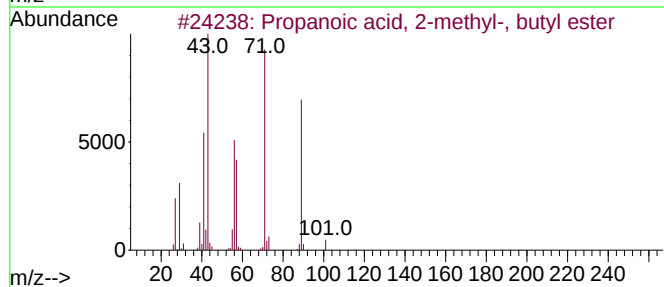
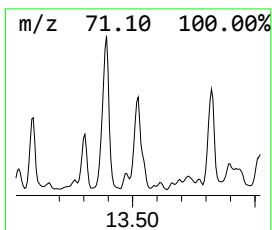
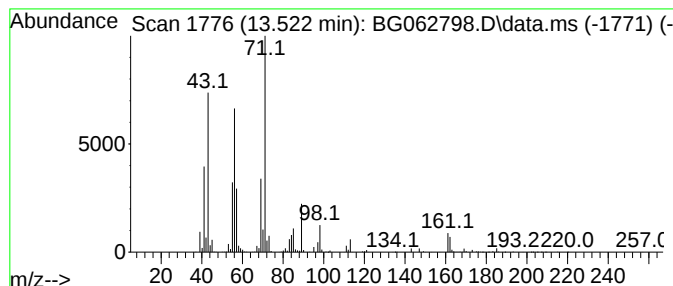
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 Propanoic acid, 2-methyl-, ... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.522	25.72 ng/ul	7578080	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	59
2			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	59
3			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	59
4			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	59
5			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
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ClientSampleId :
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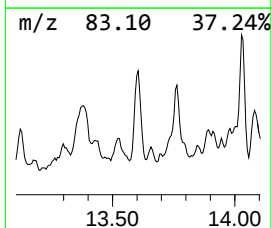
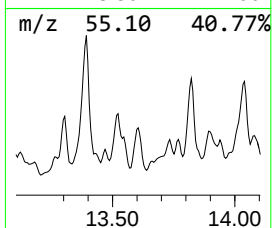
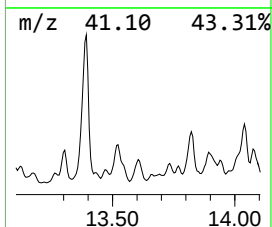
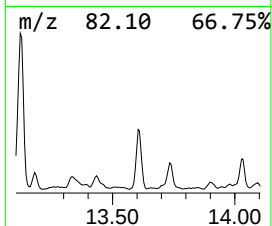
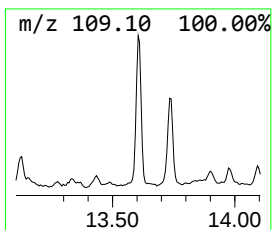
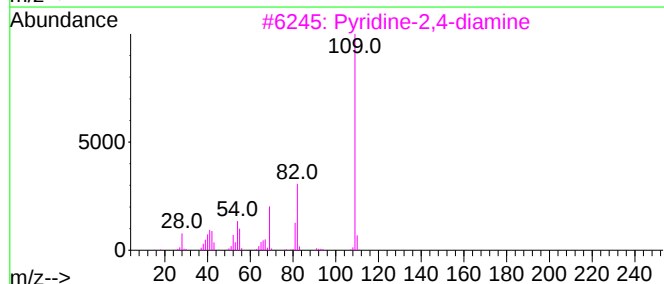
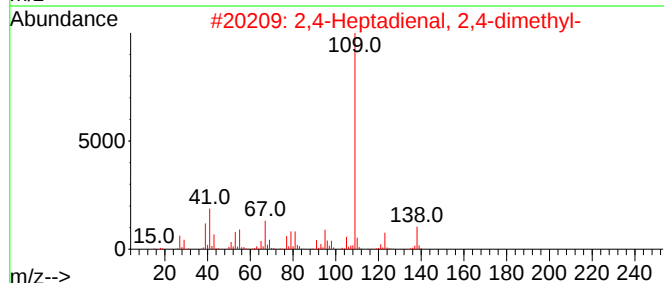
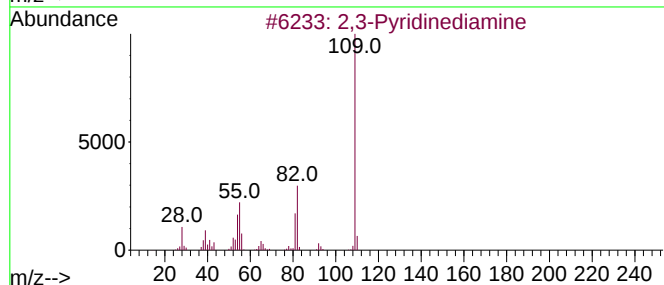
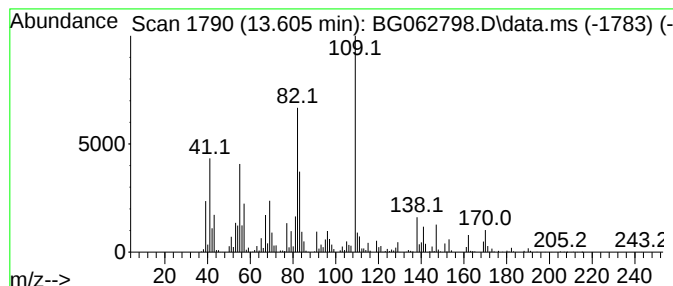
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 unknown-03 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.605	20.36 ng/ul	5999760	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3-Pyridinediamine	109	C5H7N3	000452-58-4	47
2			2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	38
3			Pyridine-2,4-diamine	109	C5H7N3	000461-88-1	38
4			Cyclopentane, (2-methylbutylidene)-	138	C10H18	053366-54-4	38
5			Metacetamol	151	C8H9NO2	000621-42-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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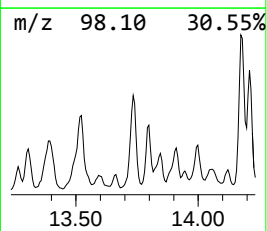
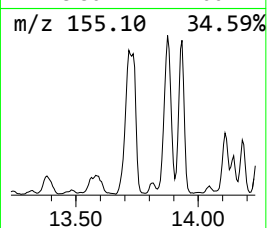
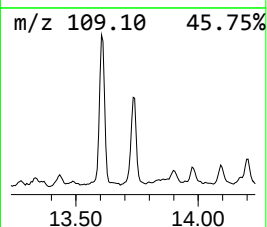
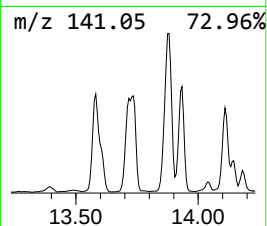
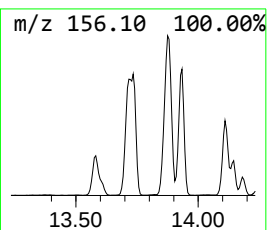
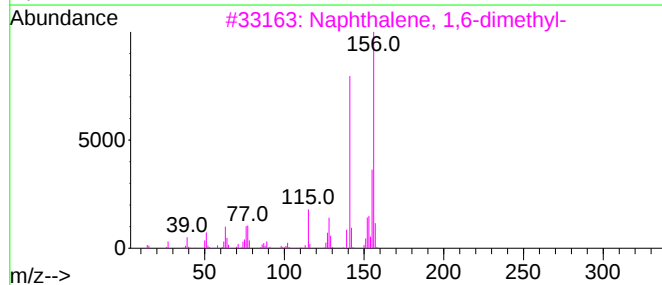
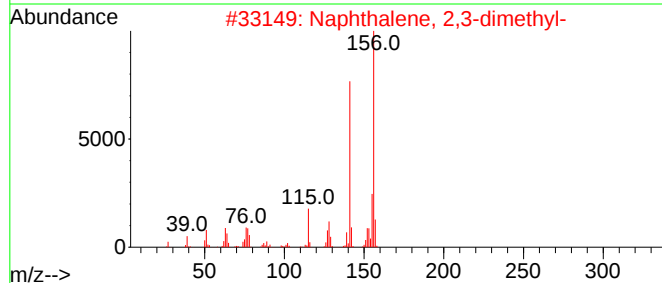
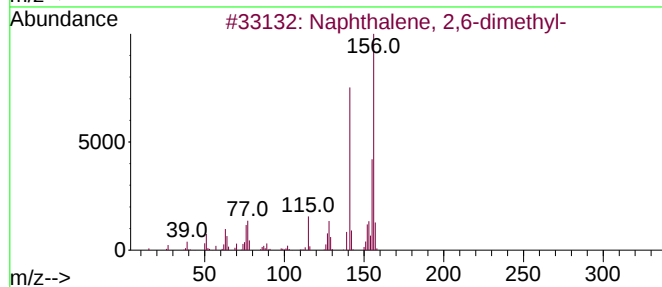
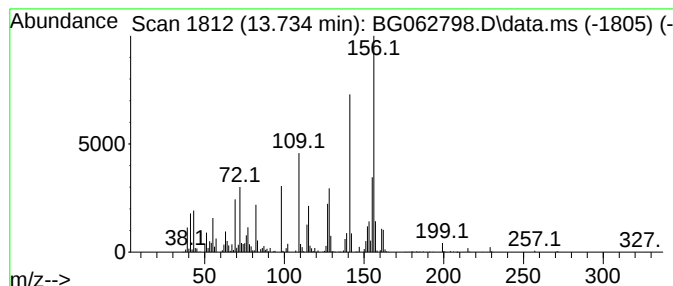
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 Naphthalene, 2,6-dimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.734	20.70 ng/ul	6098290	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	91
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	90
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	90
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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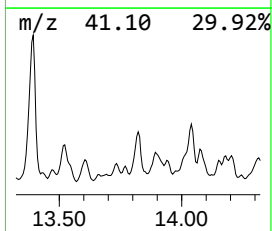
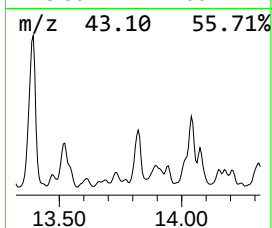
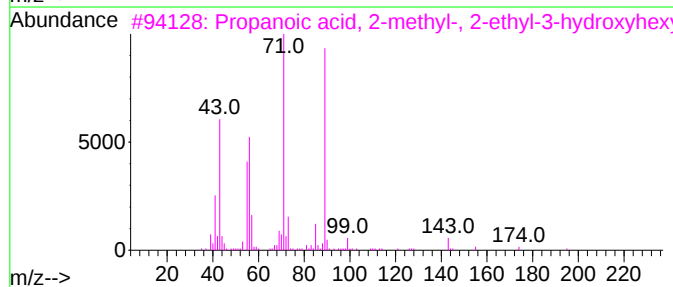
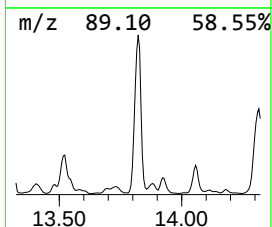
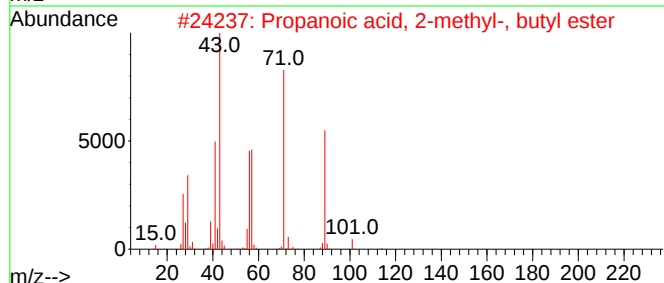
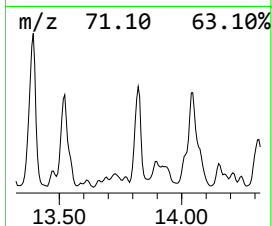
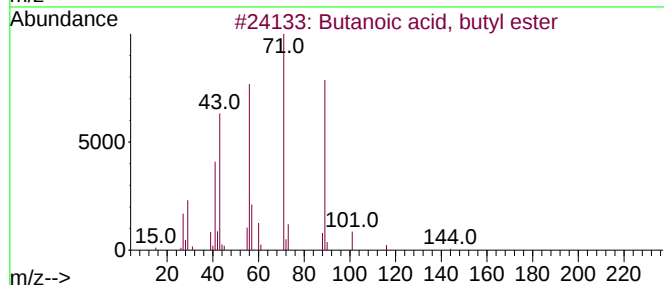
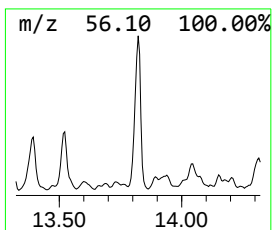
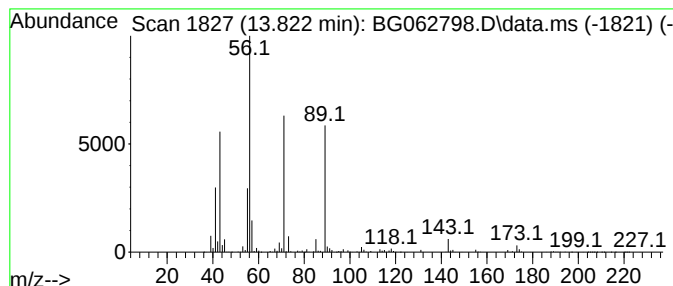
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 Butanoic acid, butyl ester Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.822	23.09 ng/ul	6804010	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	58
2			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	43
3			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	40
4			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	38
5			Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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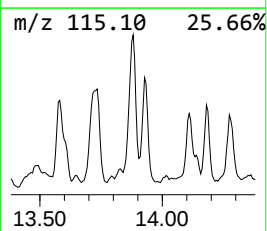
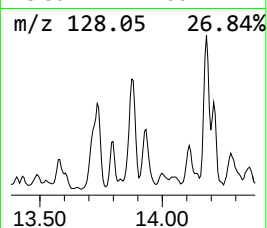
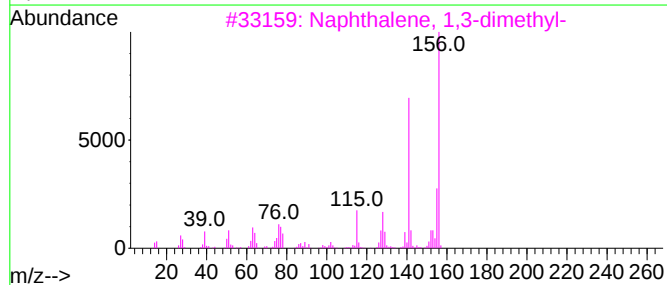
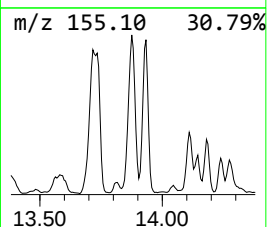
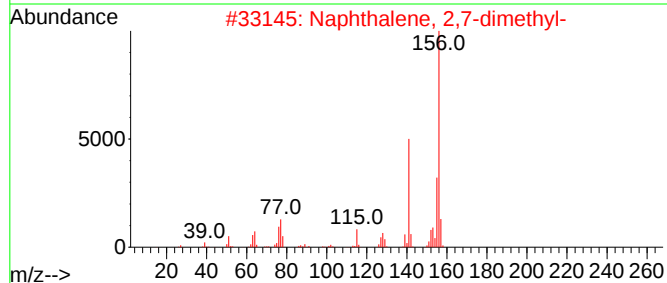
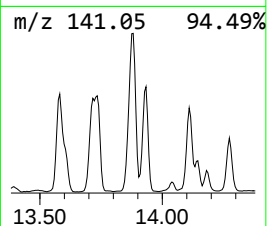
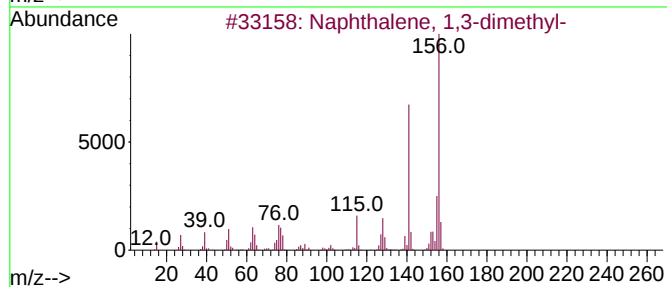
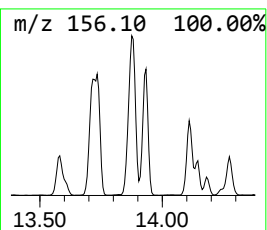
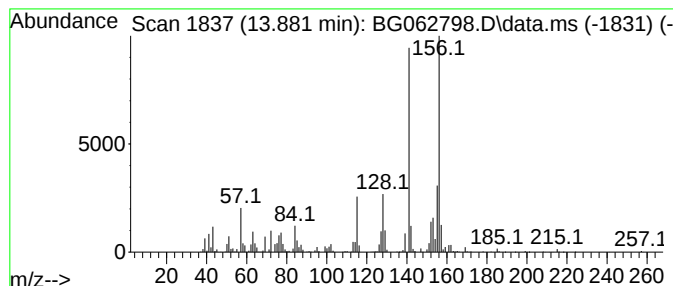
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 Naphthalene, 1,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.881	26.01 ng/ul	7665260	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

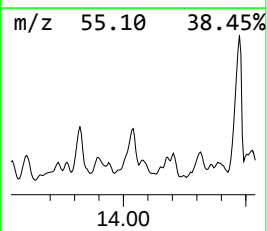
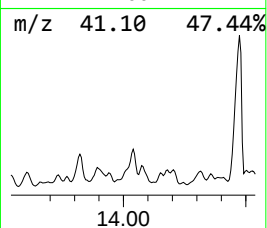
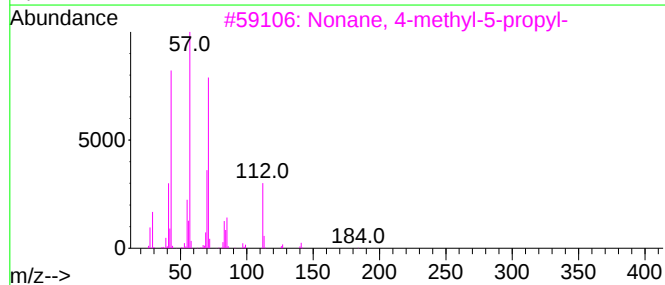
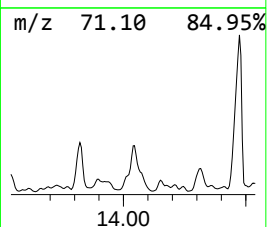
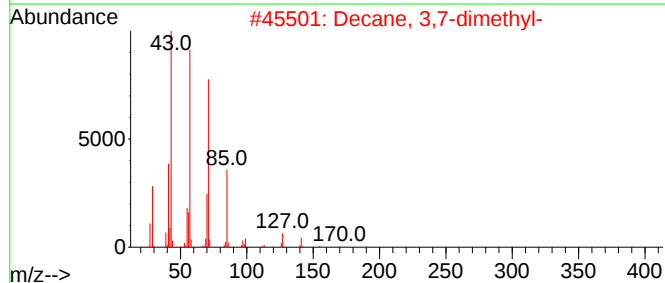
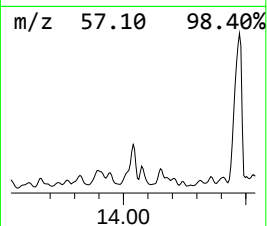
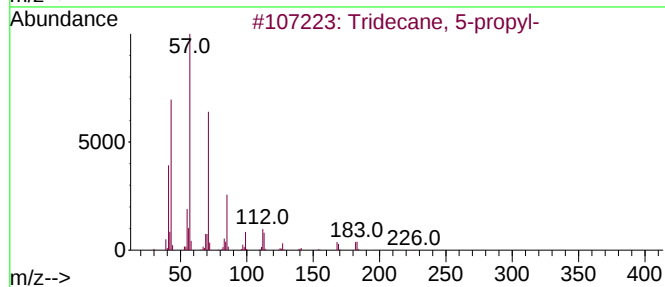
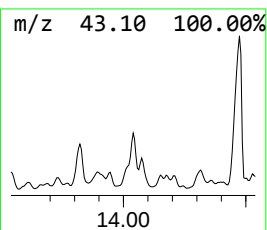
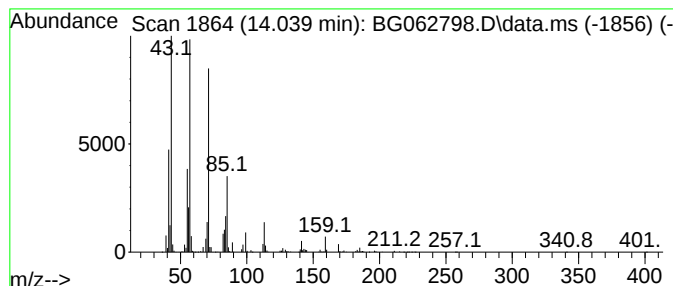
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.039	24.82 ng/ul	7312680	Acenaphthene-d10			14.551
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-		226	C16H34	055045-11-9	80
2	Decane, 3,7-dimethyl-		170	C12H26	017312-54-8	72
3	Nonane, 4-methyl-5-propyl-		184	C13H28	062185-55-1	64
4	Hexanamide, N-methallyl-		169	C10H19NO	1000340-14-0	52
5	Tridecane		184	C13H28	000629-50-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

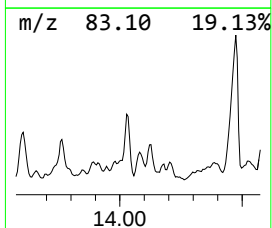
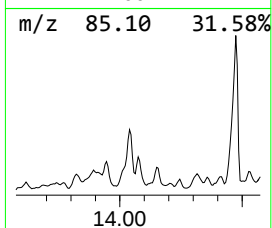
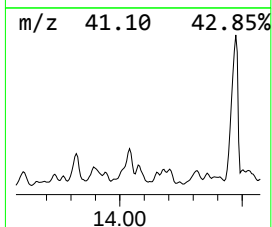
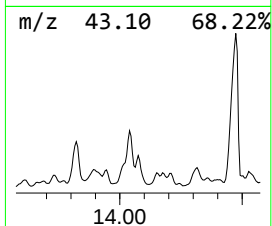
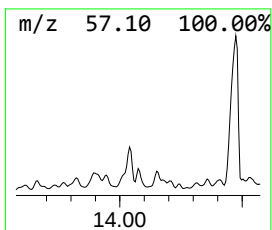
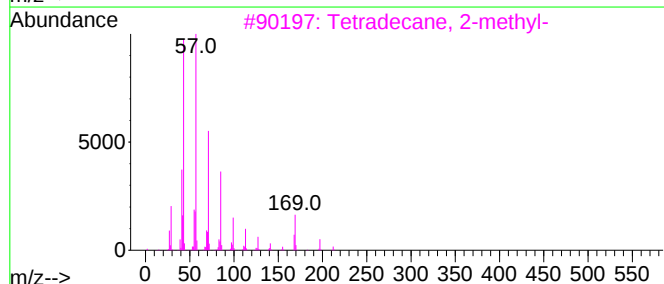
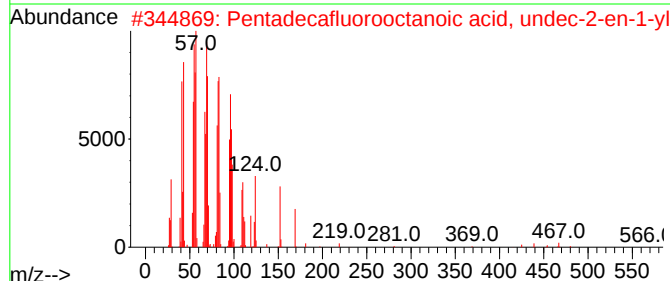
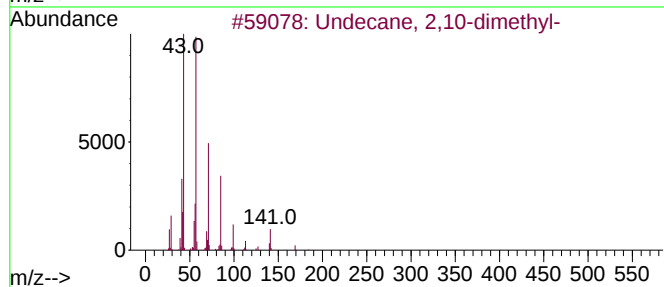
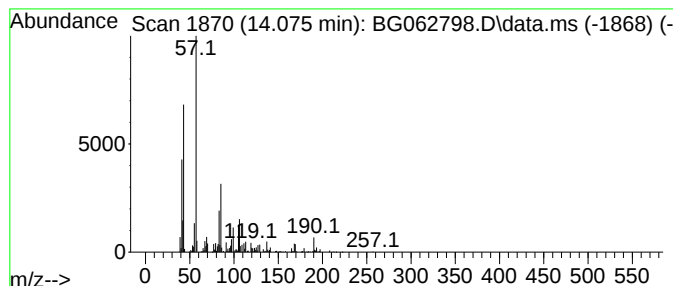
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.075	15.73 ng/ul	4636180	Acenaphthene-d10	14.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,10-dimethyl-	184	C13H28	017301-27-8	43
2	Pentadecafluorooctanoic acid, un...	566	C19H21F15O2	1000406-92-6	38
3	Tetradecane, 2-methyl-	212	C15H32	001560-95-8	38
4	Dodecane, 3-methyl-	184	C13H28	017312-57-1	35
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DDC59

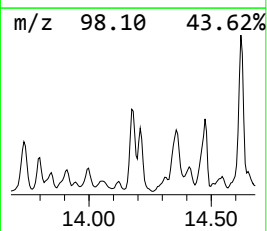
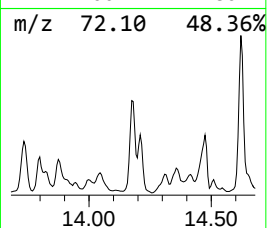
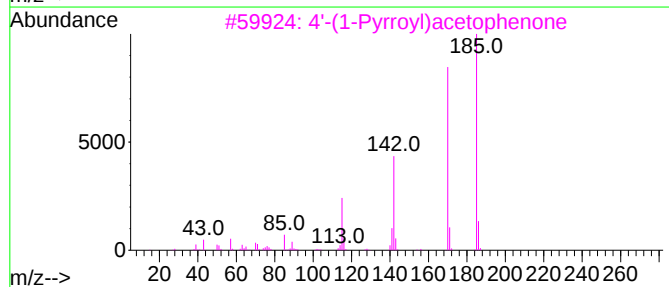
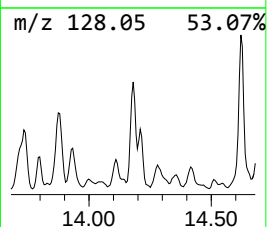
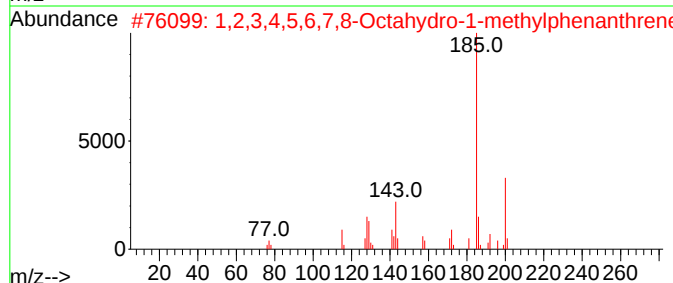
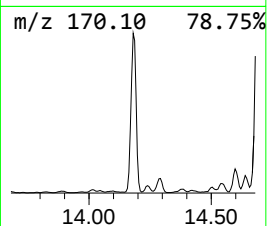
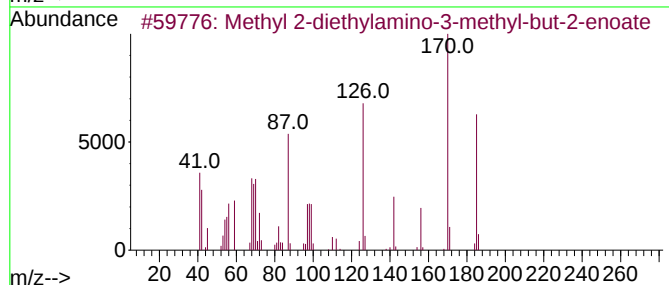
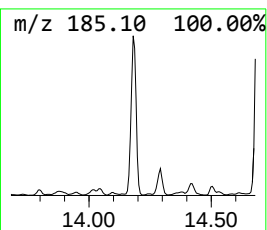
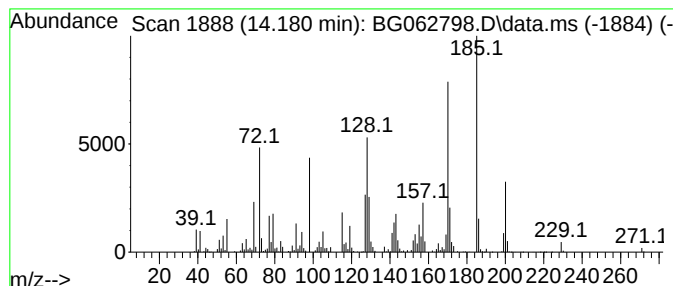
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 unknown-04 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.180	31.83 ng/ul	9379490	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	46
2			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	42
3			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	38
4			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	38
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

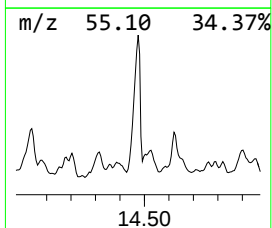
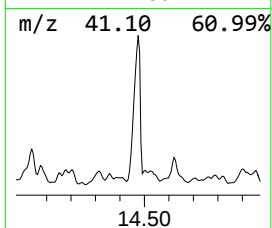
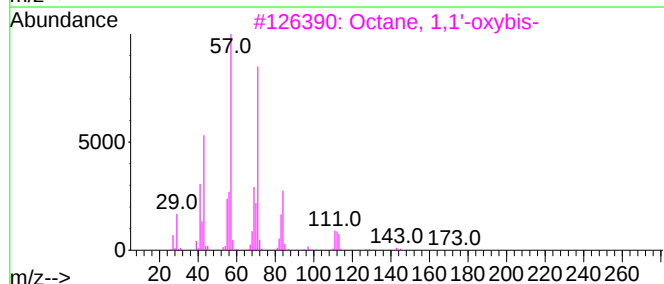
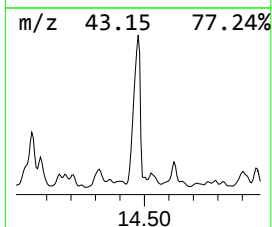
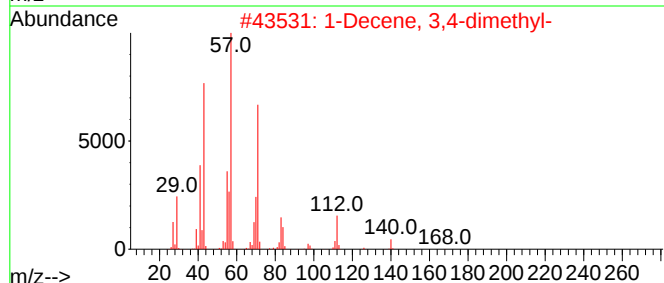
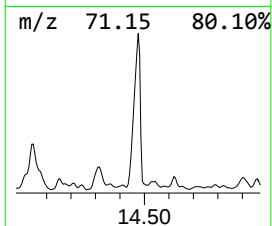
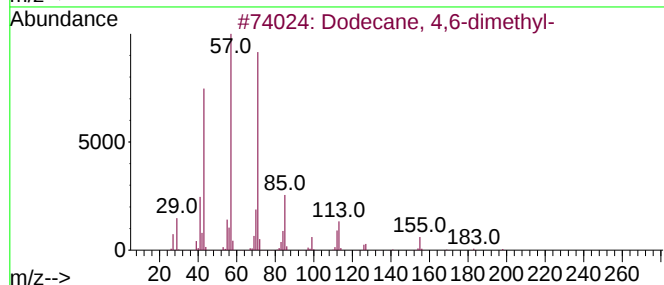
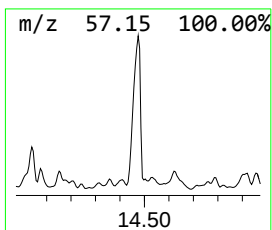
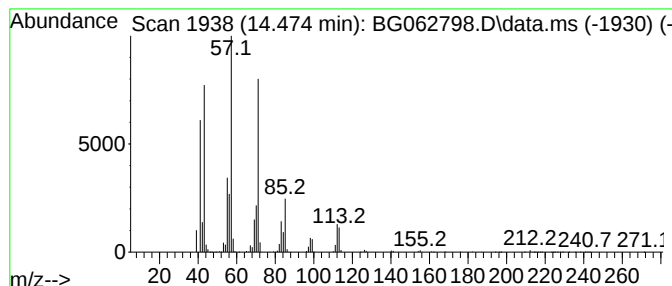
Instrument :
BNA_G
ClientSampleId :
DDC59

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.474	82.56 ng/ul	24327100	Acenaphthene-d10			14.551
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 4,6-dimethyl-		198	C14H30	061141-72-8	78
2	1-Decene, 3,4-dimethyl-		168	C12H24	050871-03-9	72
3	Octane, 1,1'-oxybis-		242	C16H34O	000629-82-3	64
4	Octane, 2,3,6-trimethyl-		156	C11H24	062016-33-5	64
5	Carbonic acid, octyl vinyl ester		200	C11H20O3	1000383-25-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

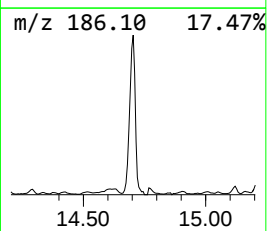
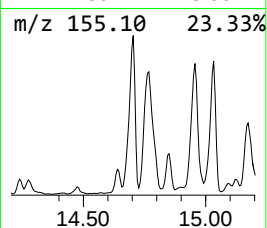
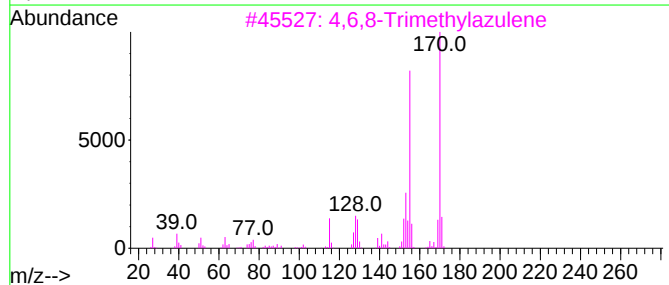
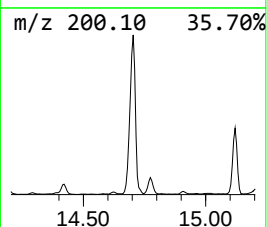
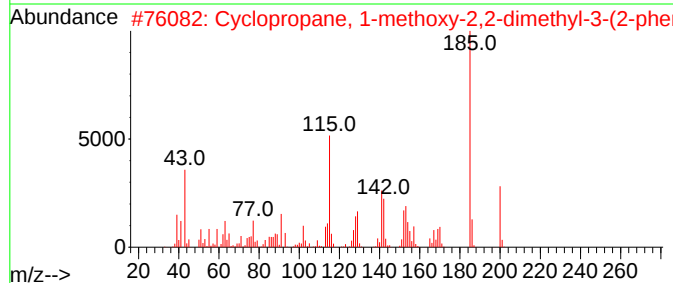
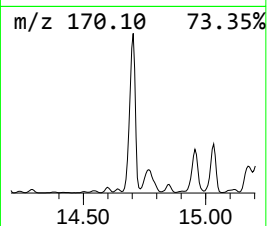
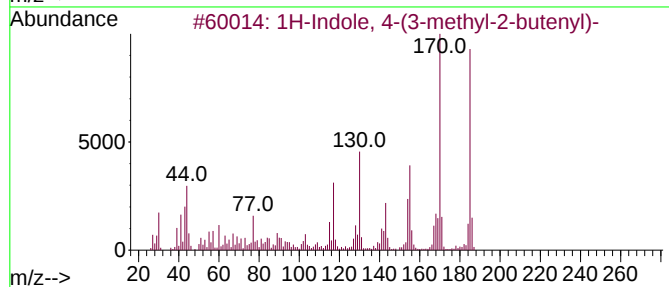
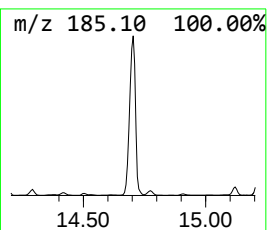
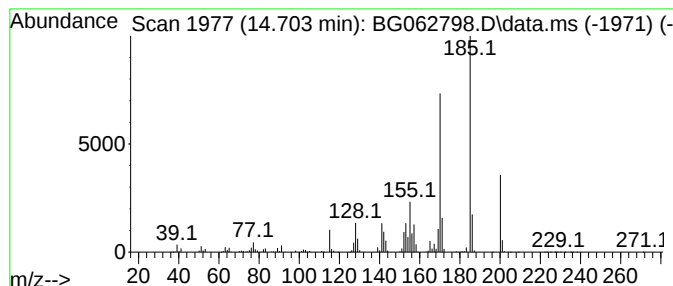
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 1H-Indole, 4-(3-methyl-2-bu... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.703	41.93 ng/ul	12355200	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indole, 4-(3-methyl-2-butenyl)-	185	C13H15N	032962-25-7	80
2			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	55
3			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	46
4			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46
5			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

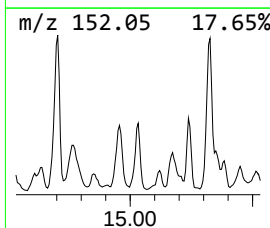
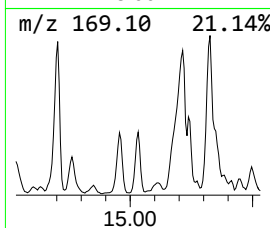
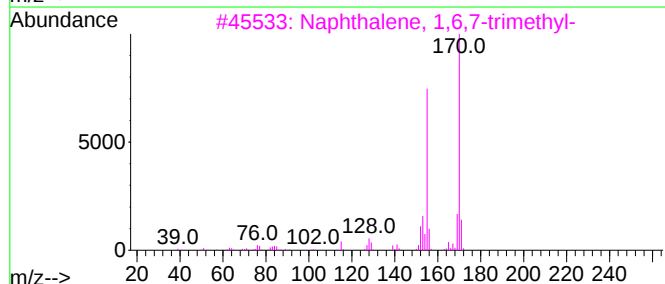
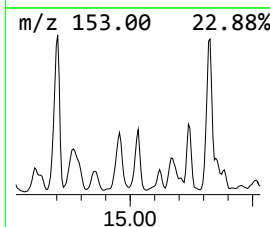
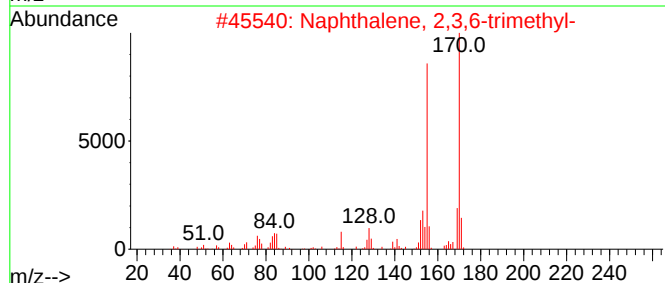
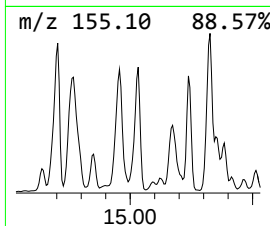
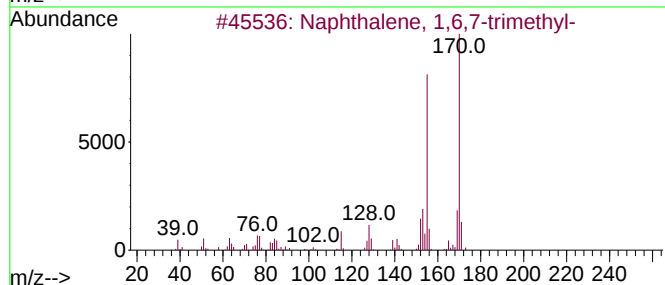
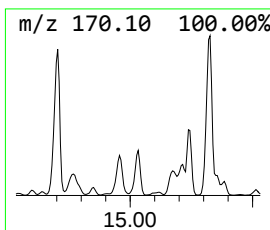
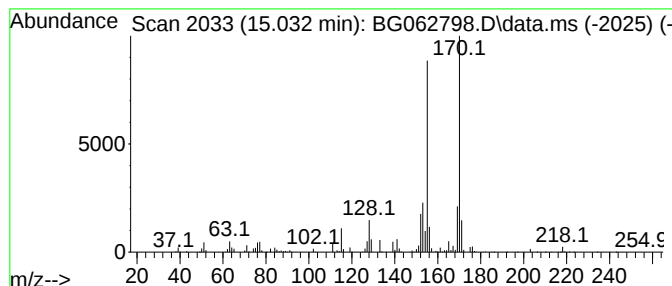
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 Naphthalene, 1,6,7-trimethyl- Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.032	12.68 ng/ul	3735100	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
5			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

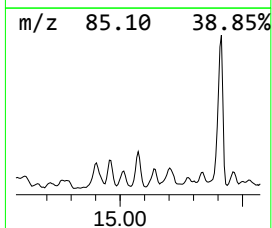
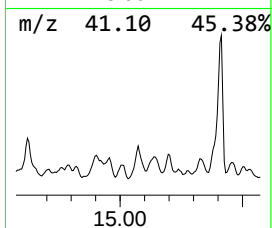
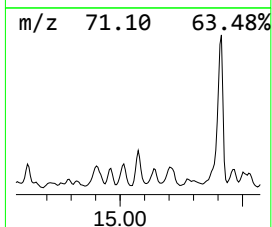
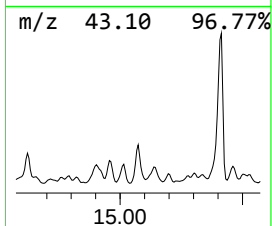
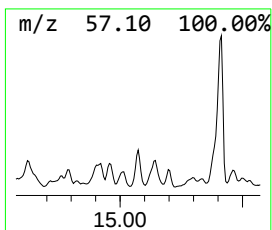
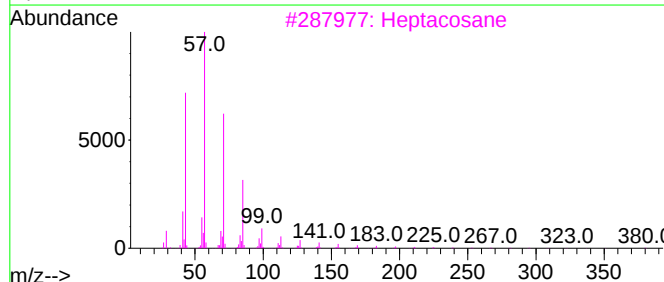
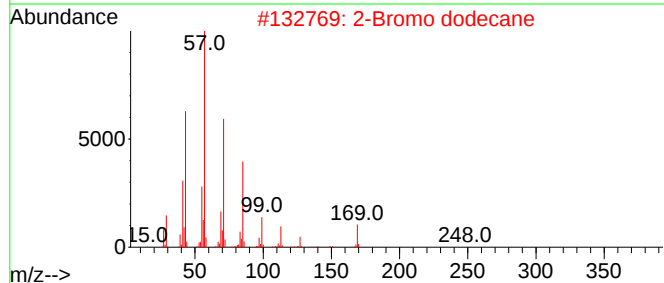
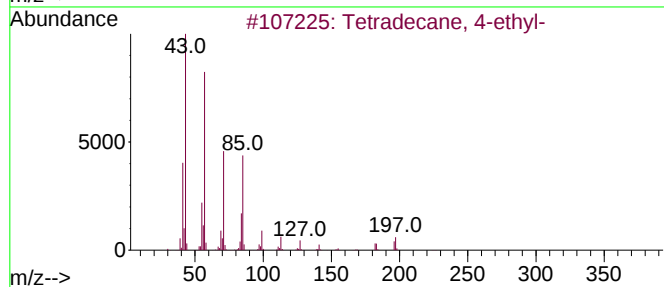
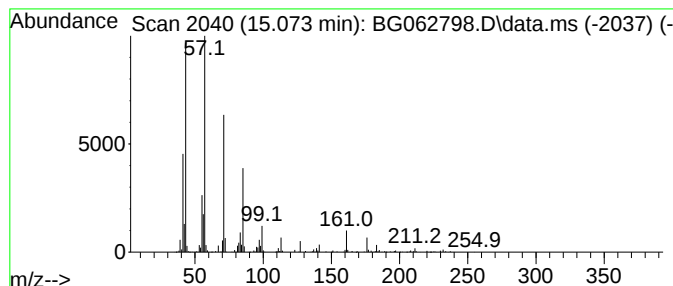
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.073	10.40 ng/ul	3064670	Acenaphthene-d10	14.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	87
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	80
3	Heptacosane	380	C27H56	000593-49-7	80
4	2-Bromotetradecane	276	C14H29Br	074036-95-6	72
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

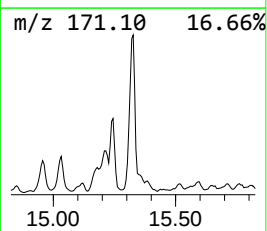
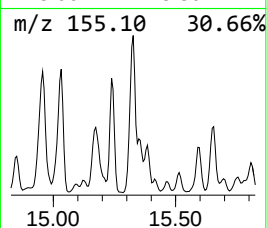
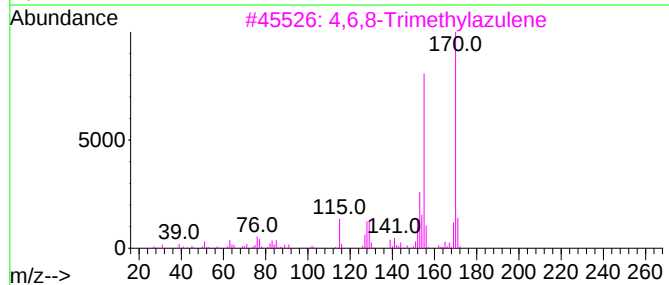
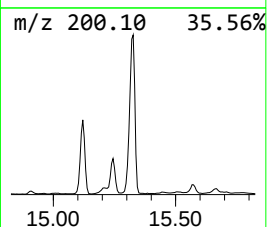
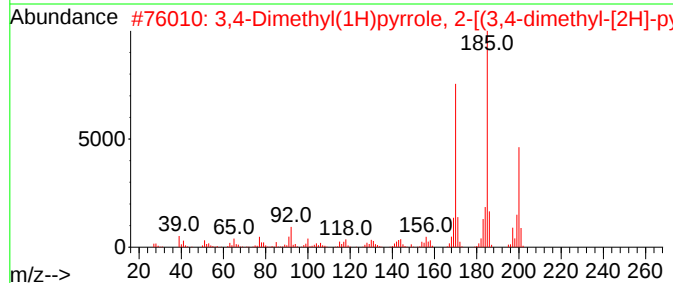
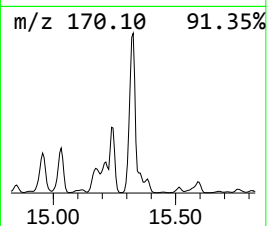
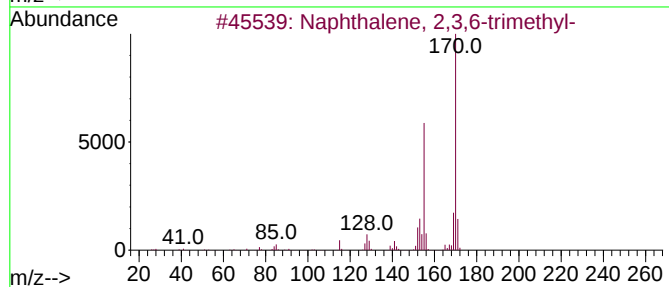
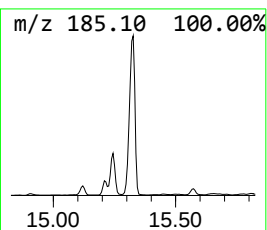
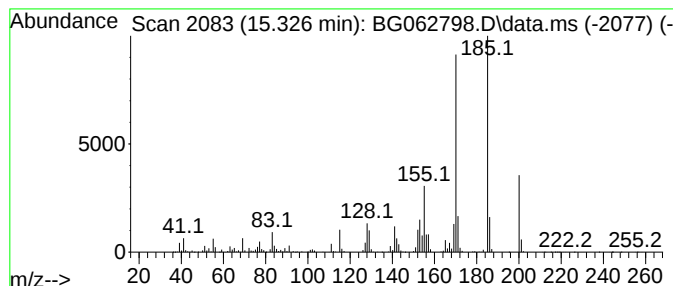
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 Naphthalene, 2,3,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.326	49.79 ng/ul	14672000	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
2			3,4-Dimethyl(1H)pyrrole, 2-[(3,4...	200	C13H16N2	065539-79-9	83
3			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

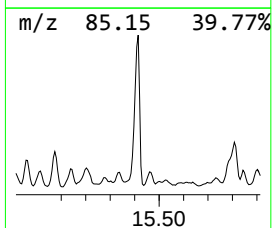
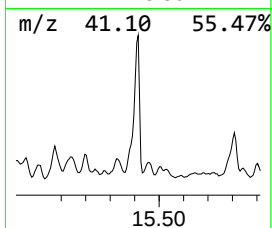
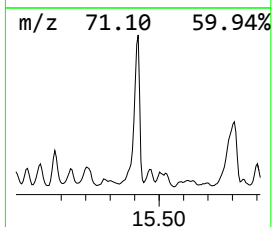
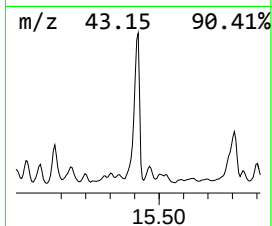
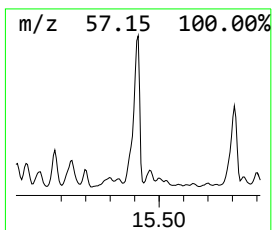
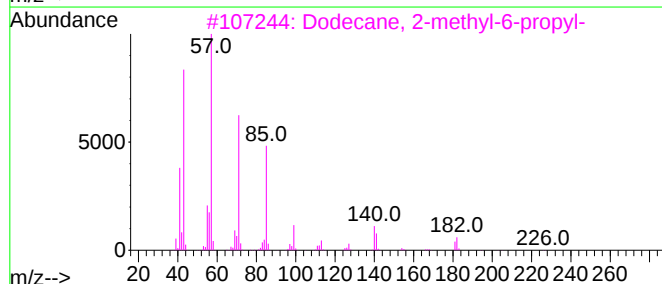
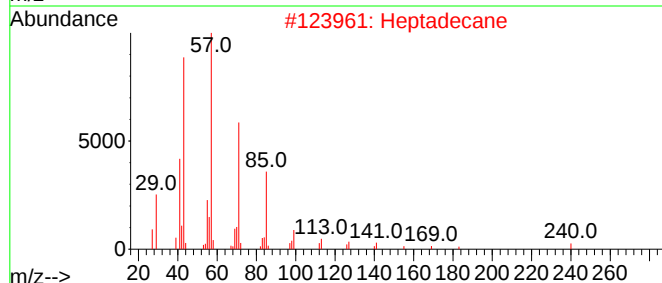
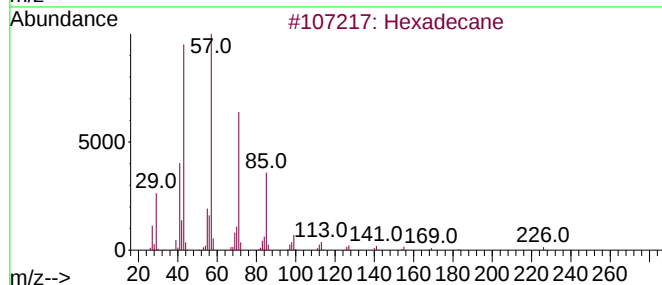
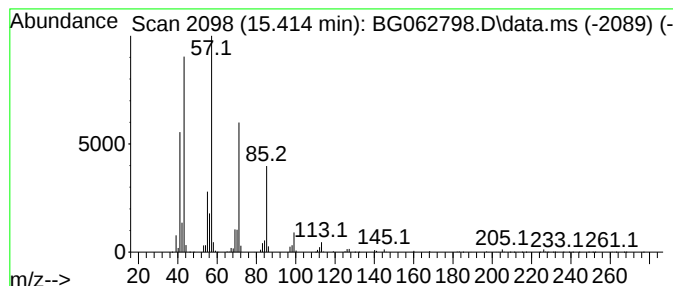
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.414	51.29 ng/ul	15114200	Acenaphthene-d10			14.551
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	93
2	Heptadecane		240	C17H36	000629-78-7	90
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	90
4	Hexadecane		226	C16H34	000544-76-3	89
5	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

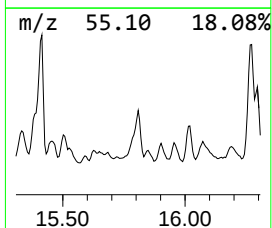
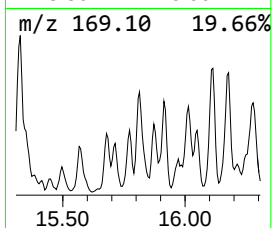
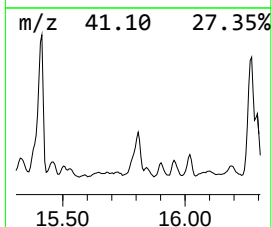
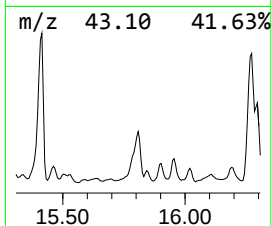
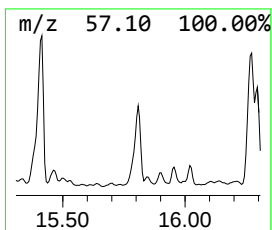
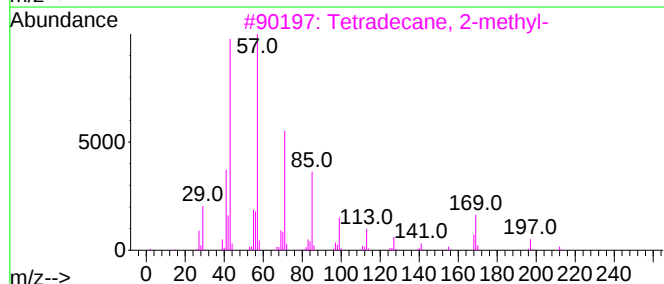
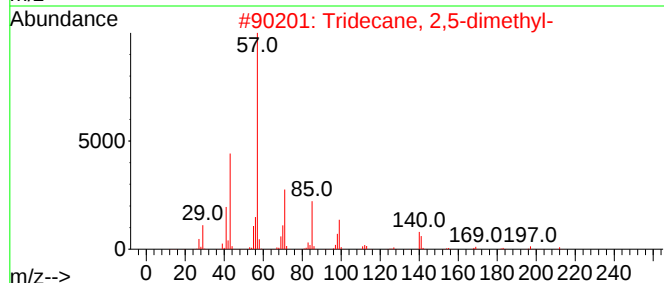
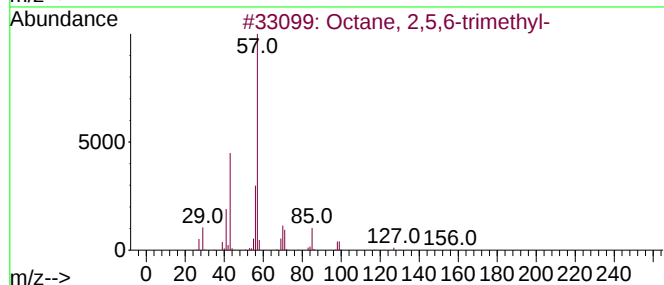
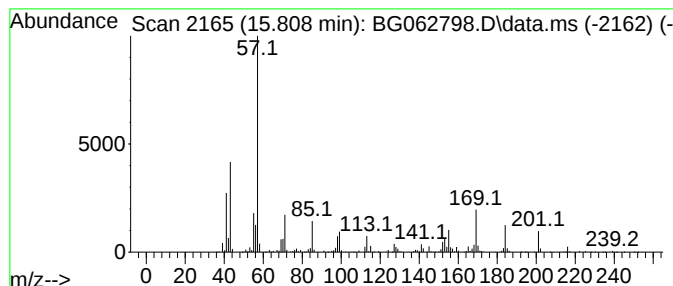
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.808	13.28 ng/ul	3914160	Acenaphthene-d10	14.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,5,6-trimethyl-	156	C11H24	062016-14-2	38
2	Tridecane, 2,5-dimethyl-	212	C15H32	056292-66-1	27
3	Tetradecane, 2-methyl-	212	C15H32	001560-95-8	22
4	Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	22
5	Heptane, 3-ethyl-	128	C9H20	015869-80-4	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

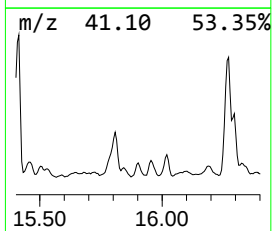
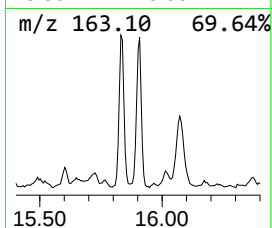
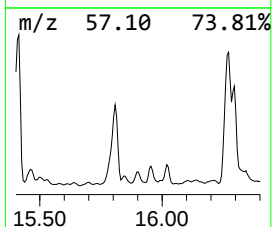
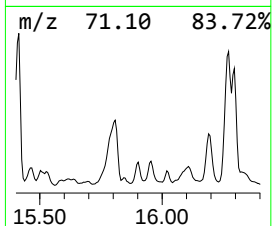
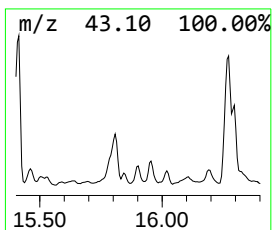
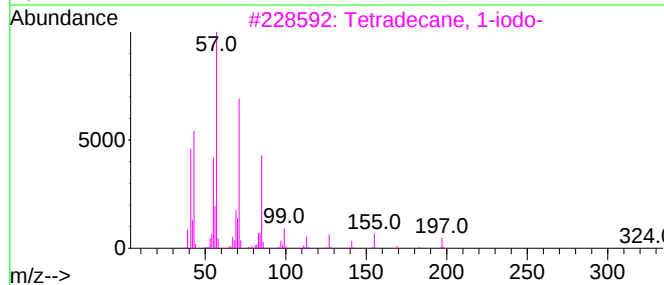
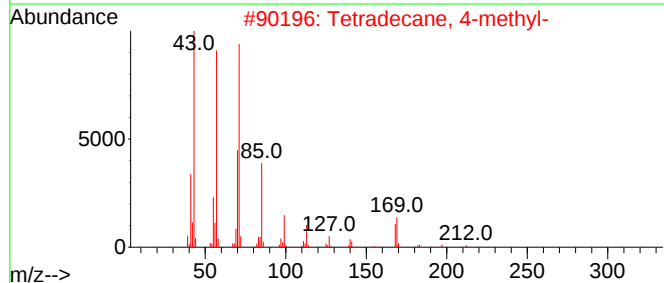
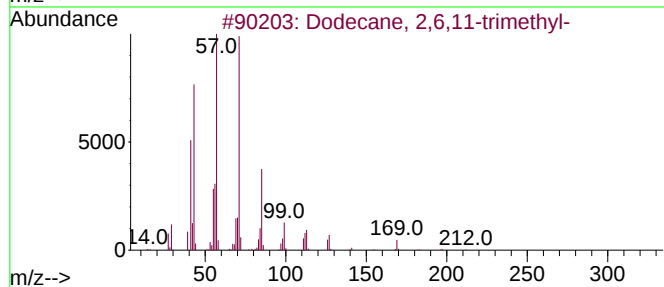
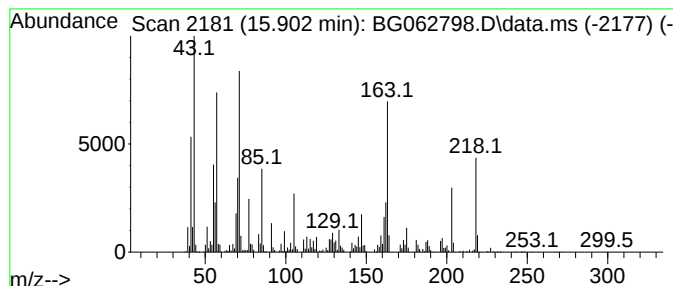
Instrument :
BNA_G
ClientSampleId :
DDC59

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.902	12.08 ng/ul	3560130	Acenaphthene-d10	14.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	46
2	Tetradecane, 4-methyl-	212	C15H32	025117-24-2	38
3	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	38
4	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	35
5	Eicosane	282	C20H42	000112-95-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

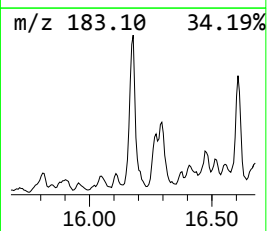
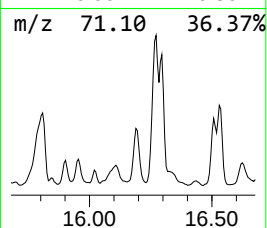
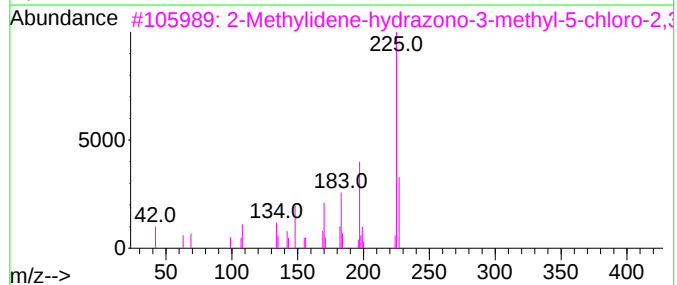
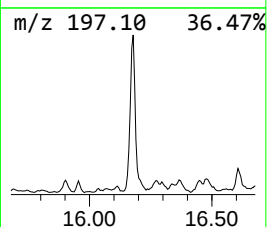
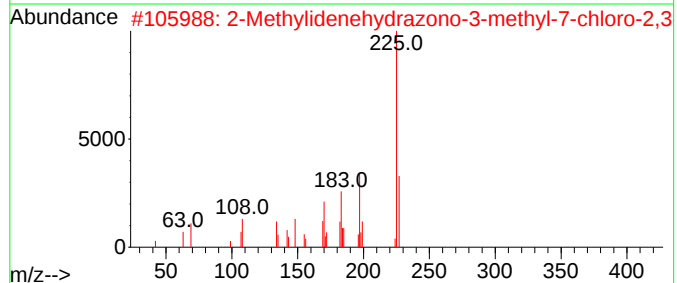
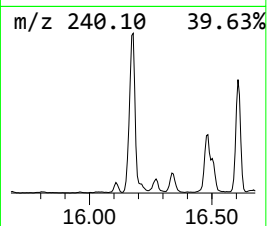
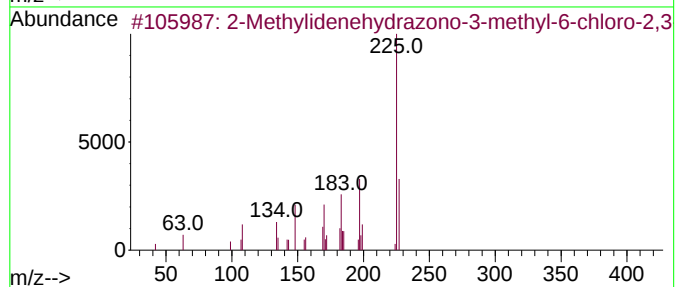
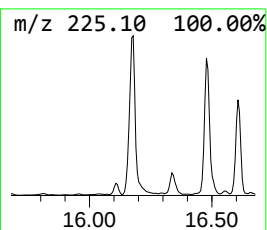
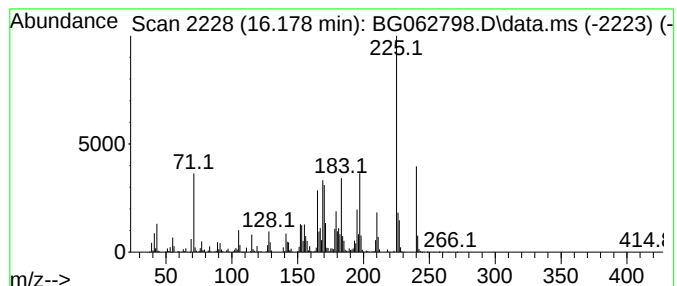
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 unknown-05 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.			
16.178	23.26 ng/ul	6385310	Phenanthrene-d10	17.341			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylidenehydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-5	46		
2	2-Methylidenehydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-6	46		
3	2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	43		
4	Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	43		
5	6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	41		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

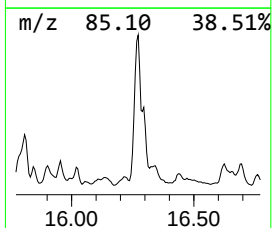
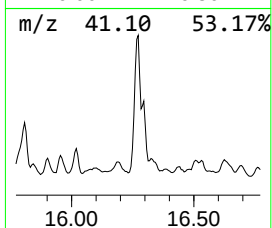
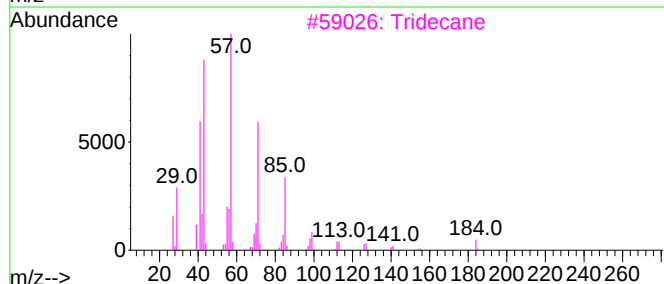
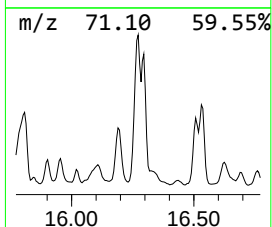
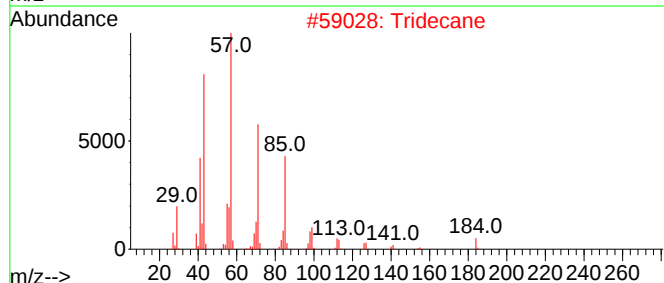
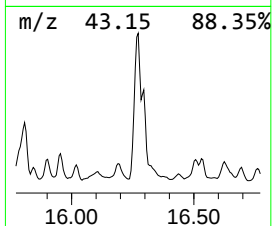
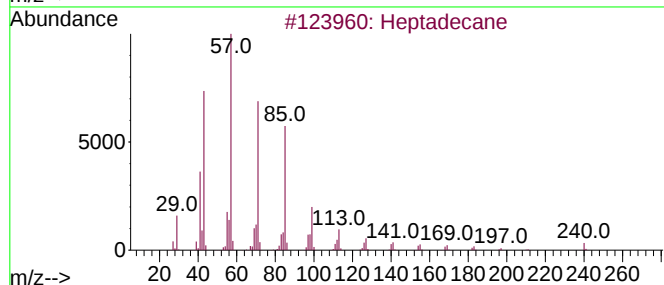
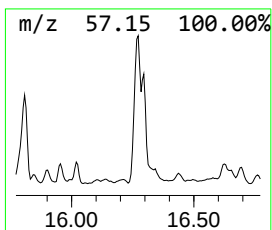
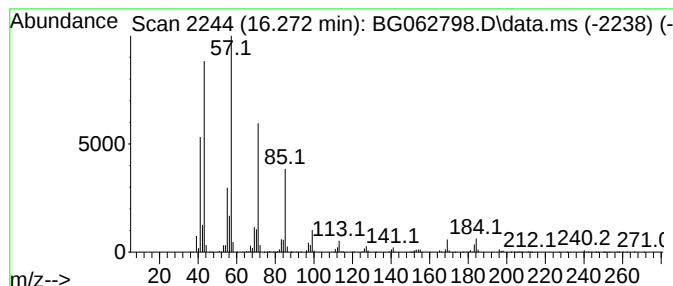
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.272	48.42 ng/ul	13291000	Phenanthrene-d10	17.341

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Tridecane	184	C13H28	000629-50-5	91
3		Tridecane	184	C13H28	000629-50-5	91
4		Nonadecane	268	C19H40	000629-92-5	86
5		Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

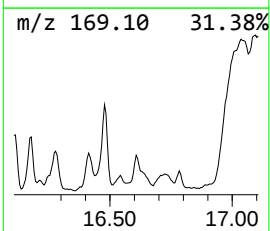
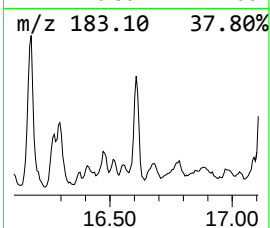
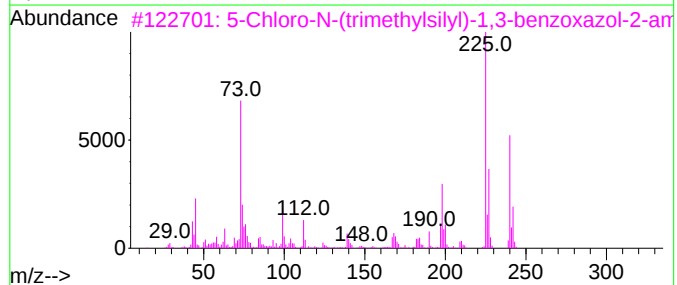
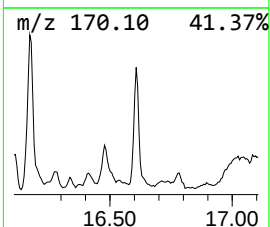
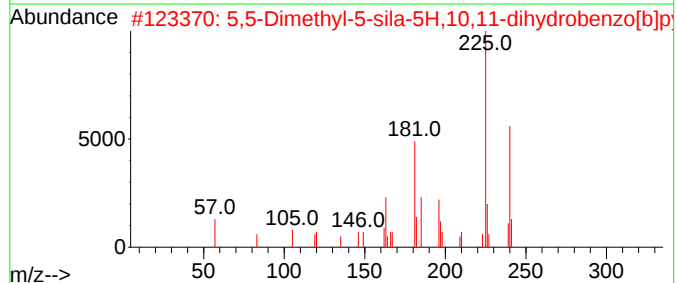
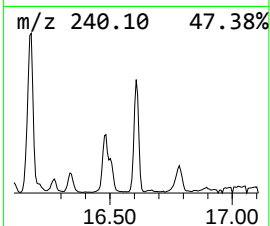
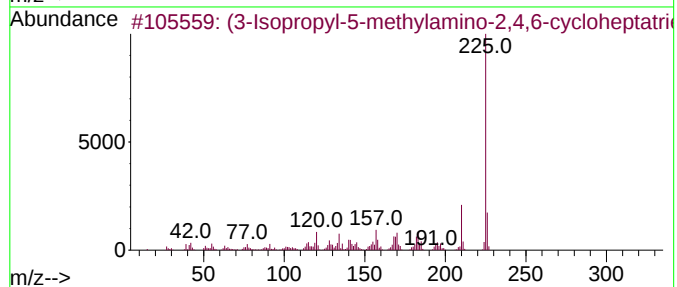
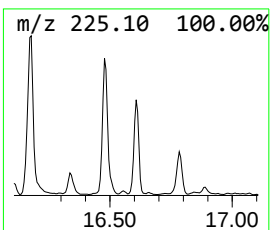
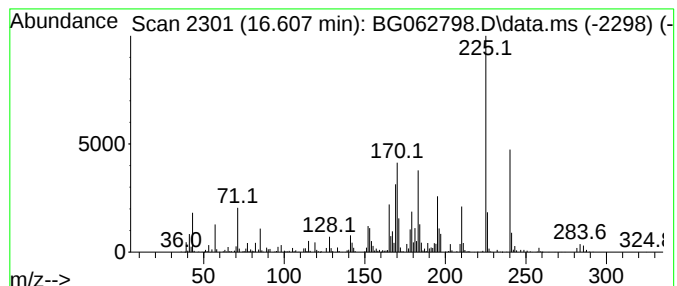
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 unknown-06 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.607	10.93 ng/ul	2999750	Phenanthrene-d10	17.341

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	38
2			5,5-Dimethyl-5-sila-5H,10,11-dih...	240	C14H16N2Si	089052-49-3	38
3			5-Chloro-N-(trimethylsilyl)-1,3-...	240	C10H13ClN2OSi	1010408-23-0	35
4			Naphthalene, 2,7-bis(1,1-dimethy...	240	C18H24	010275-58-8	35
5			4-(Dimethylamino)thiophenol, S-t...	225	C11H19NSSi	065257-43-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

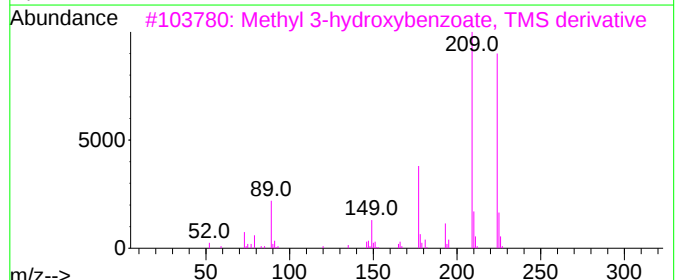
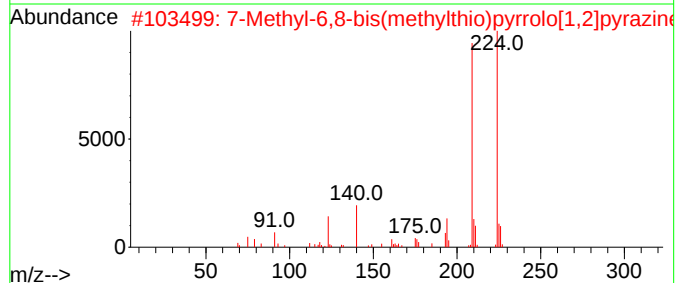
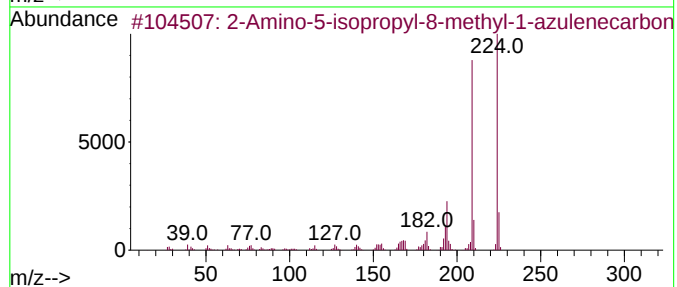
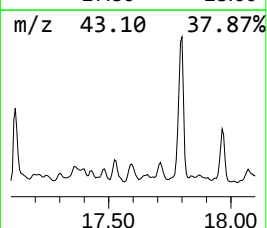
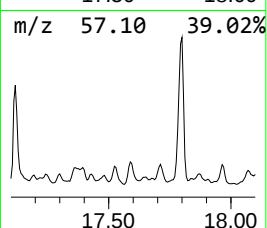
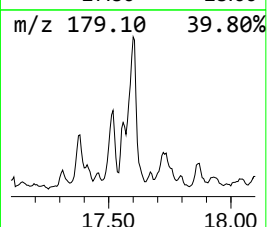
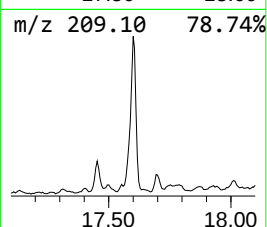
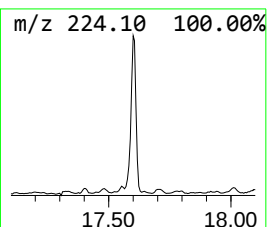
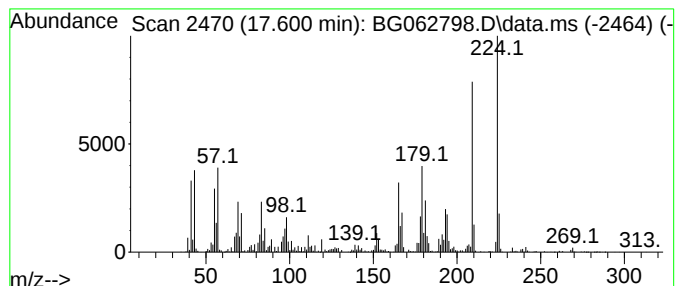
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 2-Amino-5-isopropyl-8-methy... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.600	20.00 ng/ul	5489380	Phenanthrene-d10	17.341	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	86
2	7-Methyl-6,8-bis(methylthio)pyrr...	224	C10H12N2S2	084201-40-1	52
3	Methyl 3-hydroxybenzoate, TMS de...	224	C11H16O3Si	027798-50-1	52
4	Carbazole, 2,4,7-trimethyl-	209	C15H15N	078787-89-0	46
5	1,5,6,7-Tetramethyl-3-phenylbicy...	224	C17H20	126584-00-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

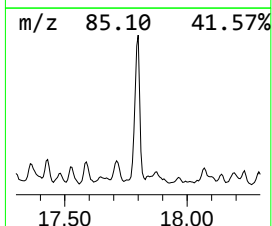
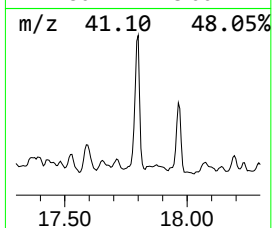
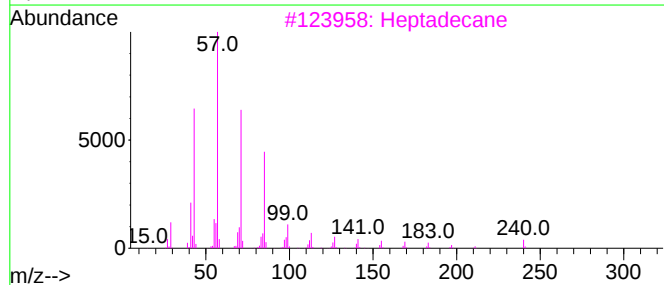
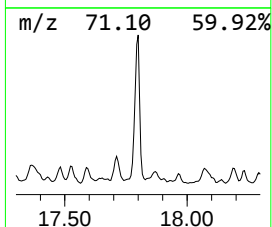
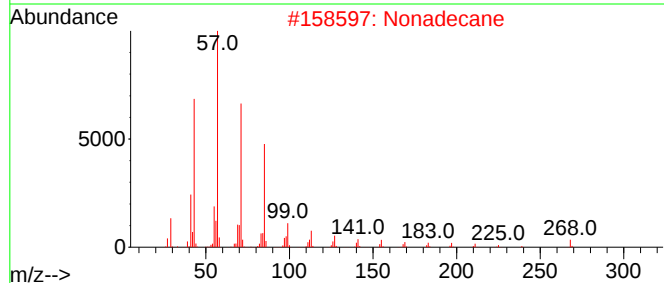
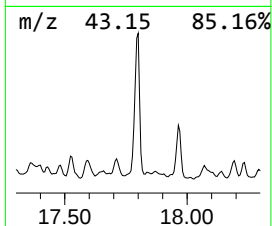
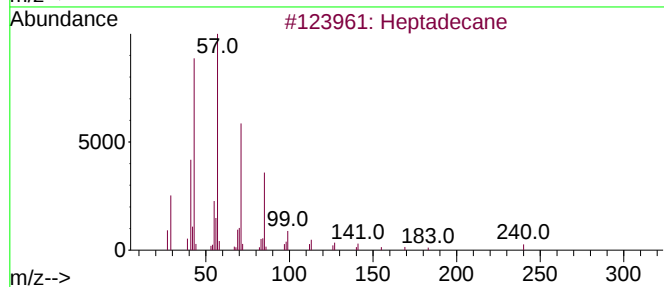
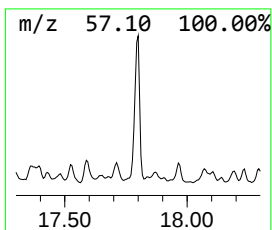
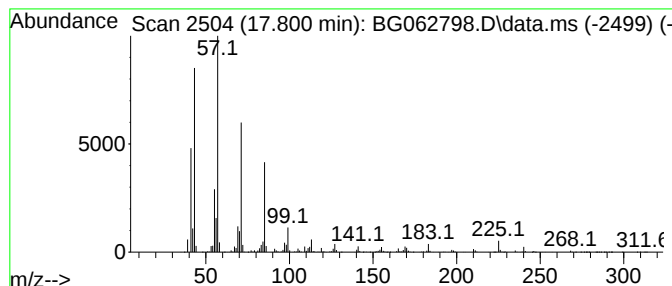
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.800	34.58 ng/ul	9492230	Phenanthrene-d10		17.341	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Nonadecane		268	C19H40	000629-92-5	95
3	Heptadecane		240	C17H36	000629-78-7	95
4	Heptadecane		240	C17H36	000629-78-7	95
5	Hexadecane		226	C16H34	000544-76-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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Acq On : 13 Sep 2024 21:36
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Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

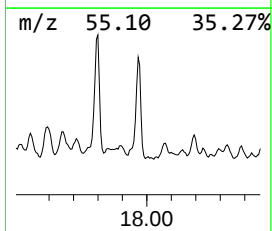
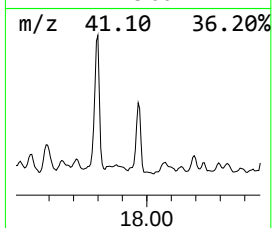
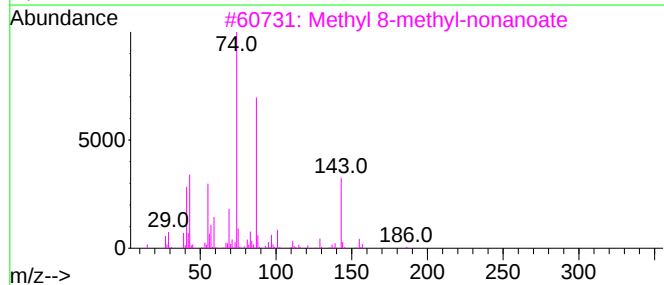
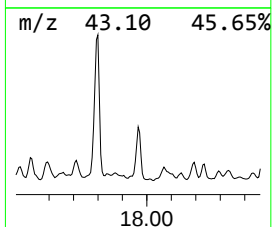
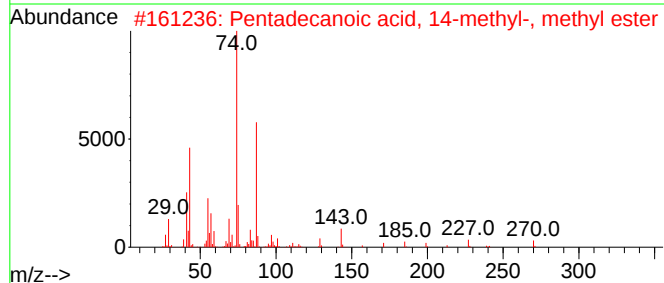
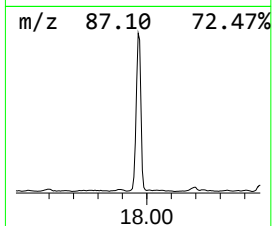
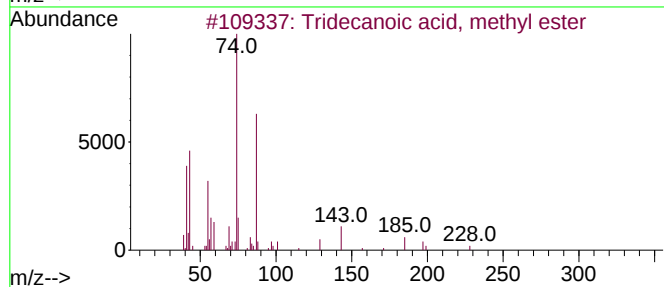
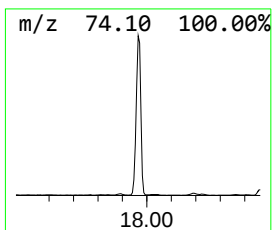
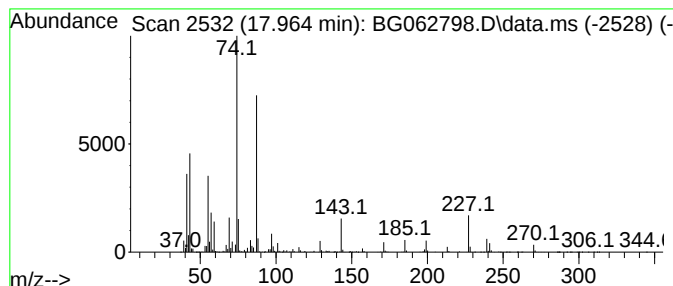
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 Tridecanoic acid, methyl ester Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.964	21.60 ng/ul	5927850	Phenanthrene-d10	17.341	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	95
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
3	Methyl 8-methyl-nonanoate	186	C11H22O2	1000452-00-9	89
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5	Methyl tetradecanoate	242	C15H30O2	000124-10-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

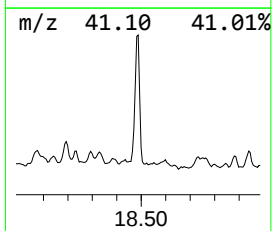
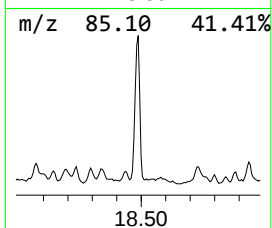
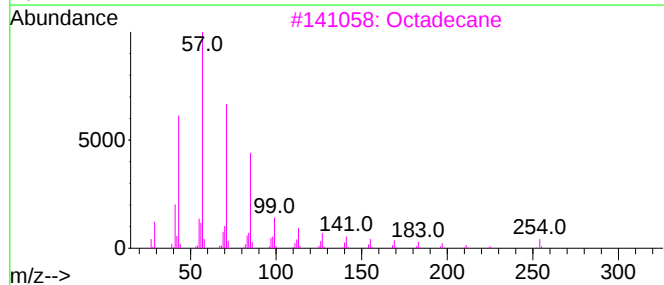
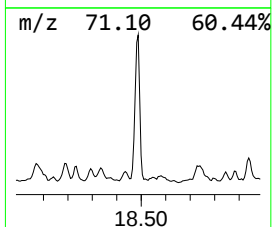
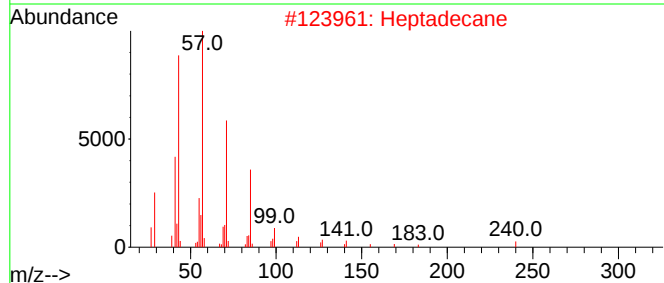
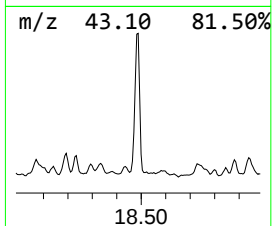
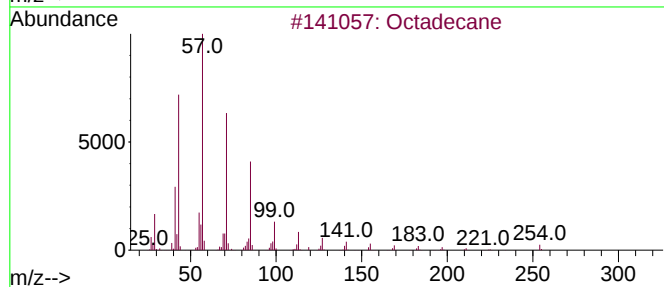
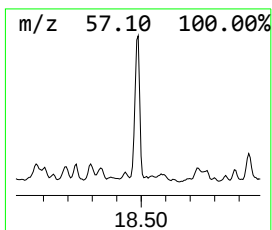
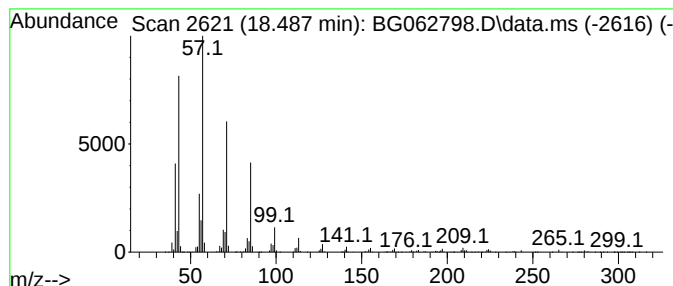
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.487	23.38 ng/ul	6415770	Phenanthrene-d10		17.341	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	95
2	Heptadecane		240	C17H36	000629-78-7	95
3	Octadecane		254	C18H38	000593-45-3	93
4	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	93
5	1-Iodoundecane		282	C11H23I	004282-44-4	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

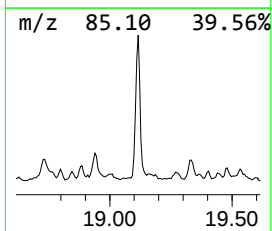
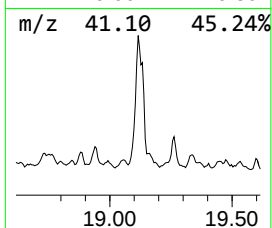
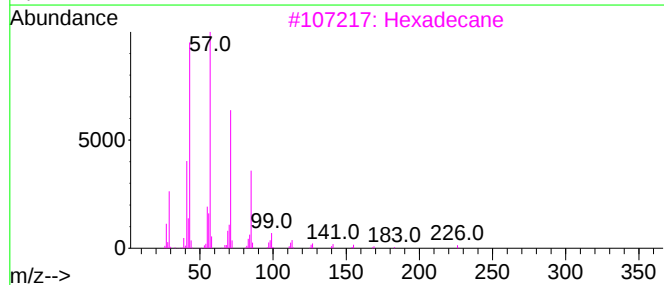
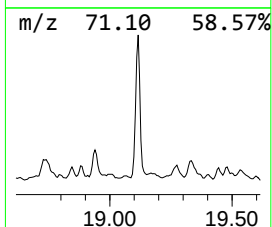
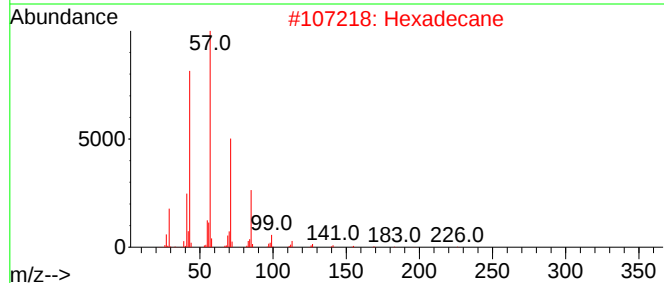
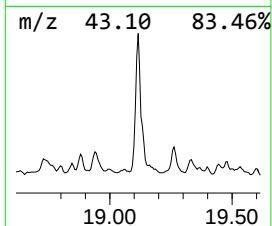
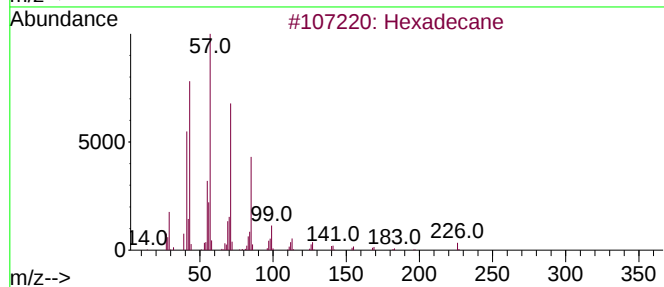
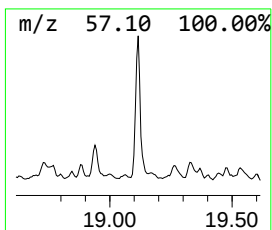
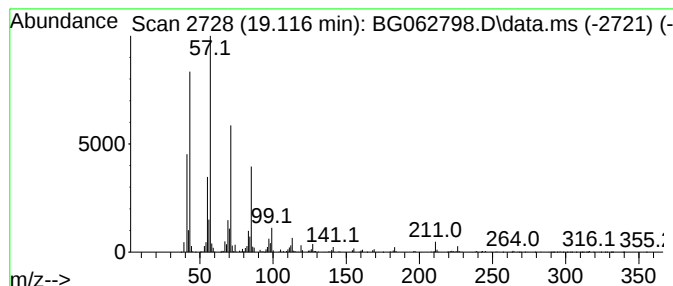
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.116	29.29 ng/ul	8038340	Phenanthrene-d10		17.341	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	98
2	Hexadecane		226	C16H34	000544-76-3	95
3	Hexadecane		226	C16H34	000544-76-3	93
4	Nonadecane		268	C19H40	000629-92-5	91
5	Nonadecane		268	C19H40	000629-92-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

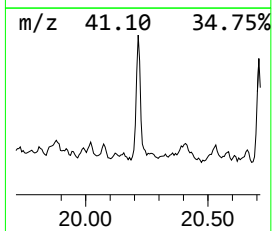
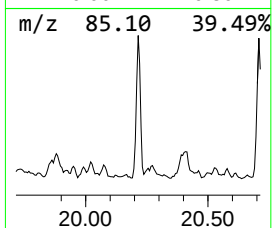
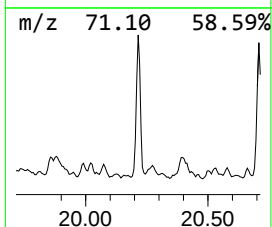
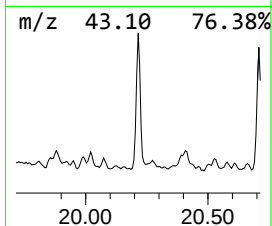
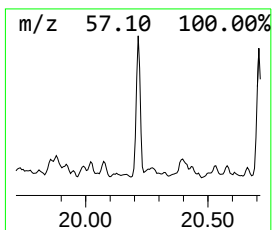
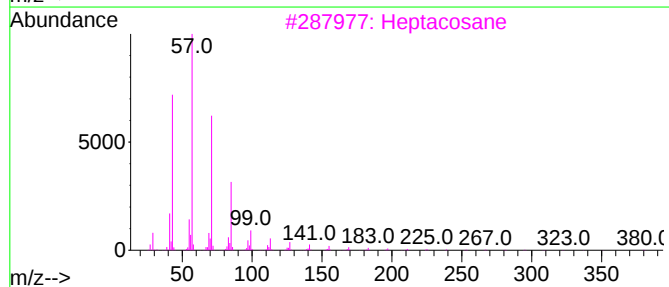
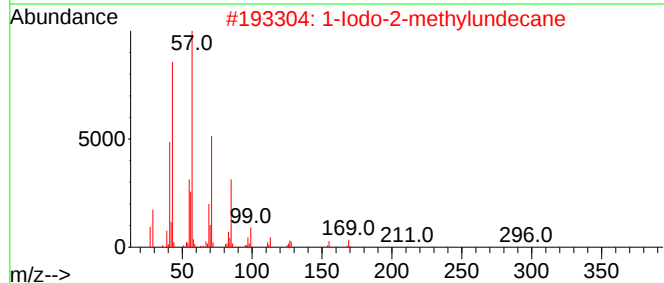
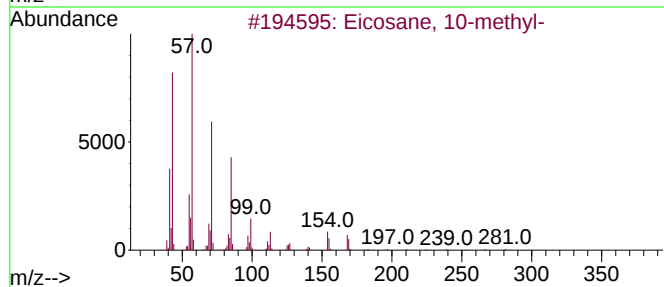
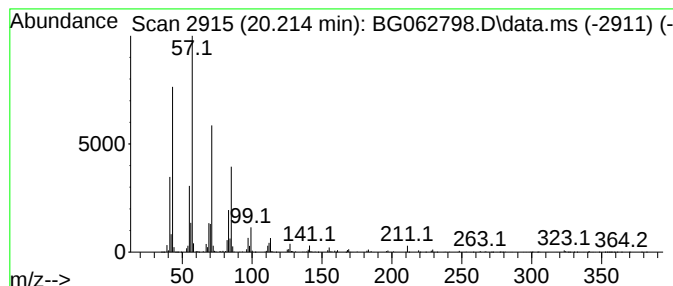
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.215	23.30 ng/ul	3840190	Chrysene-d12	21.531

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
3		Heptacosane	380	C27H56	000593-49-7	87
4		Hexacosane	366	C26H54	000630-01-3	86
5		Nonadecane	268	C19H40	000629-92-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

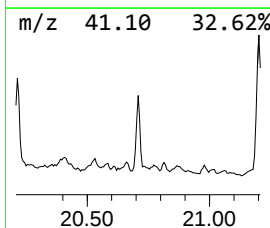
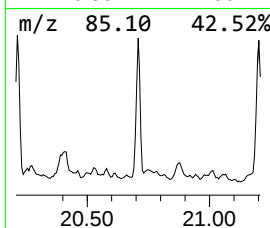
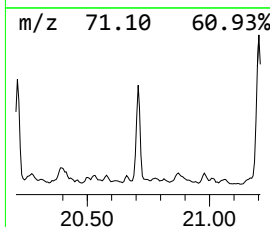
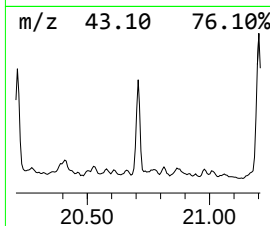
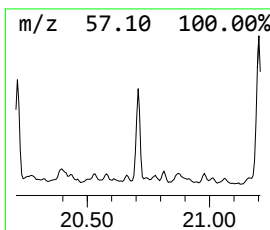
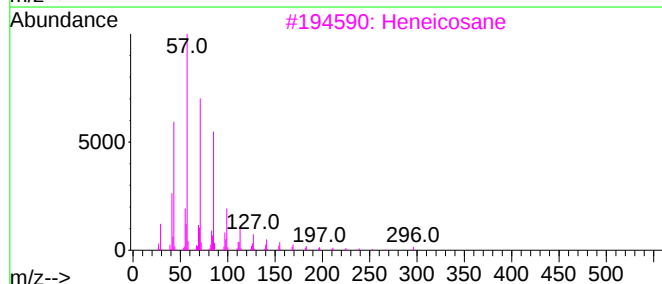
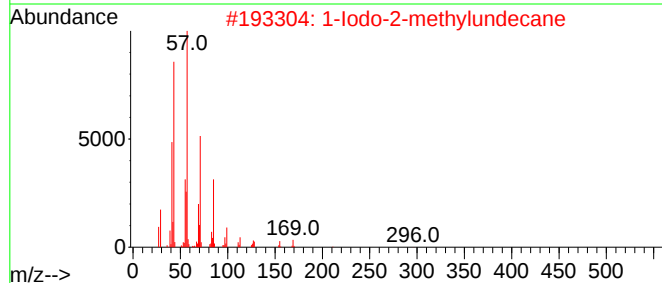
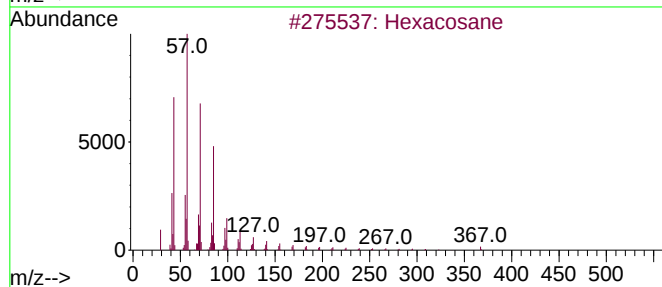
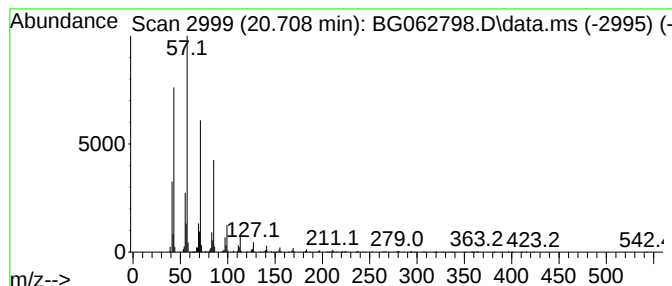
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.708	17.61 ng/ul	2901600	Chrysene-d12	21.531	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
3	Heneicosane	296	C21H44	000629-94-7	90
4	Octacosane	394	C28H58	000630-02-4	90
5	Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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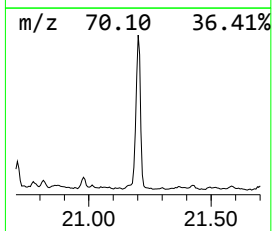
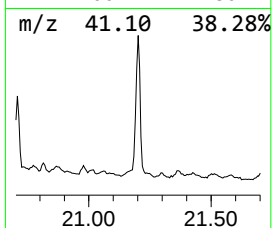
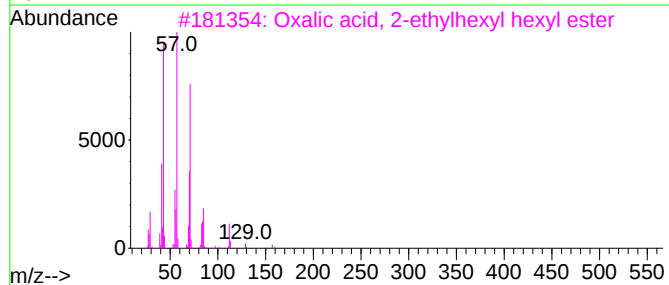
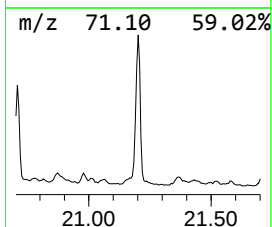
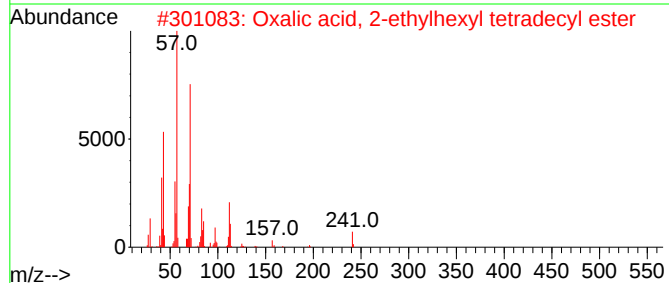
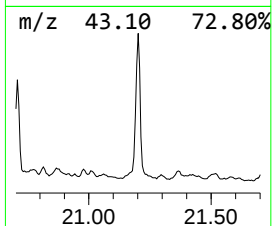
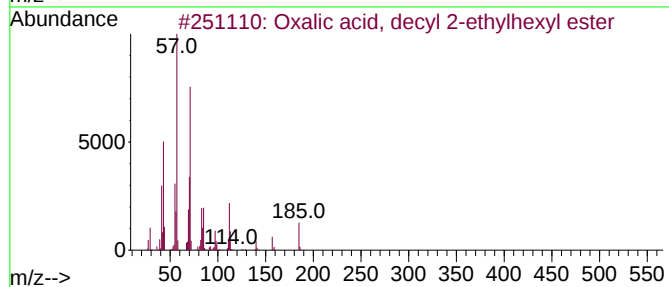
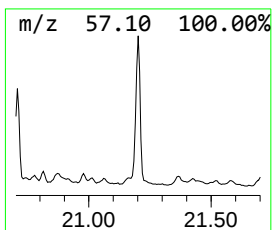
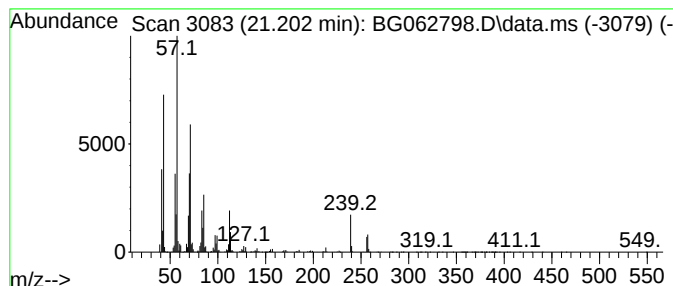
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 Oxalic acid, decyl 2-ethylh... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.202	44.38 ng/ul	7313450	Chrysene-d12	21.531

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	80
2			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	80
3			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	52
4			Tetratetracontane	619	C44H90	007098-22-8	49
5			Hexadecane	226	C16H34	000544-76-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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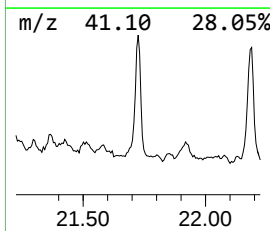
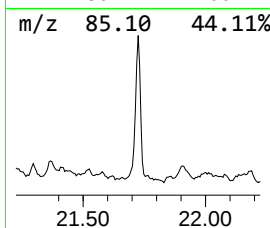
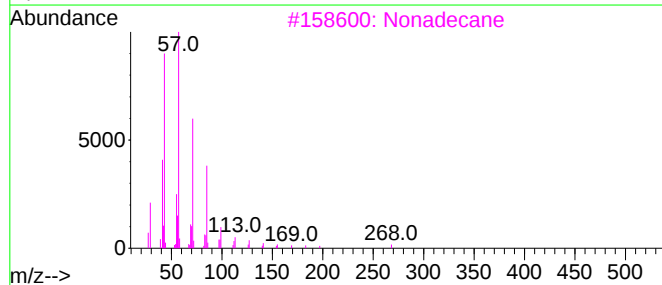
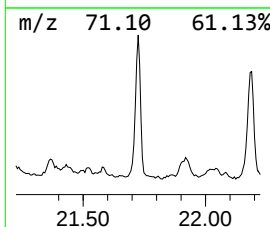
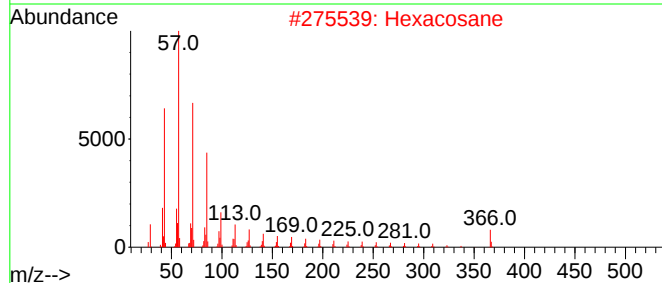
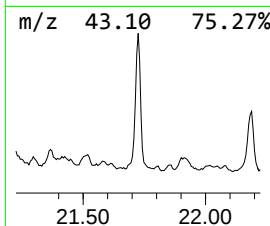
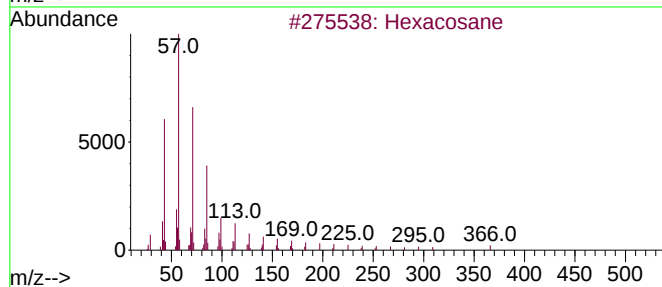
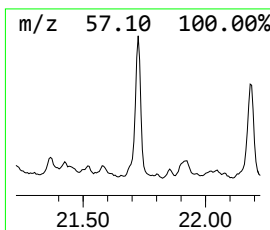
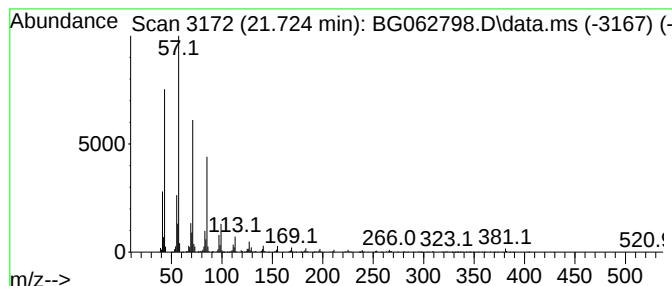
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.724	20.00 ng/ul	3296200	Chrysene-d12	21.531

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	97
2			Hexacosane	366	C26H54	000630-01-3	95
3			Nonadecane	268	C19H40	000629-92-5	95
4			Heneicosane	296	C21H44	000629-94-7	95
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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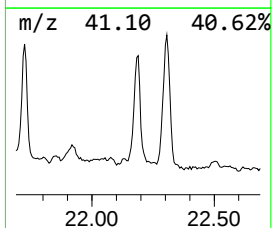
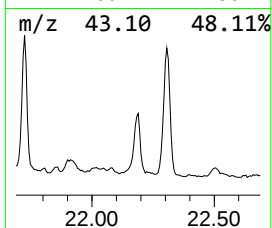
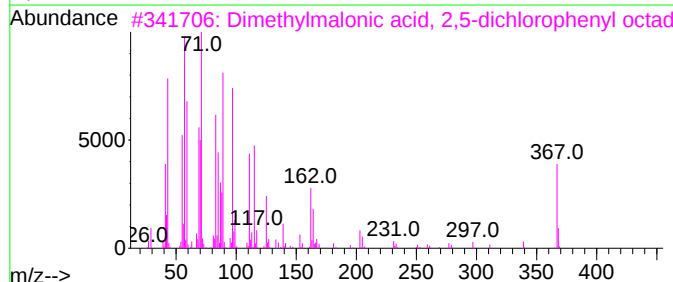
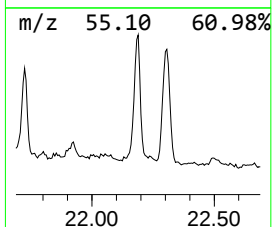
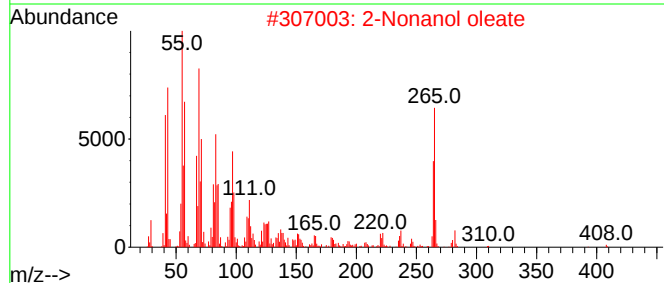
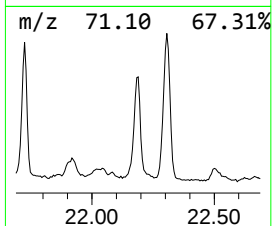
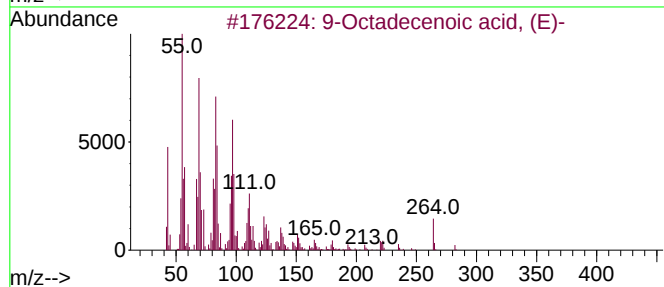
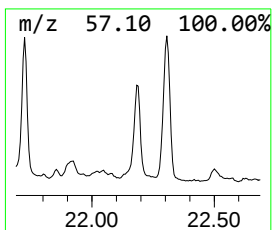
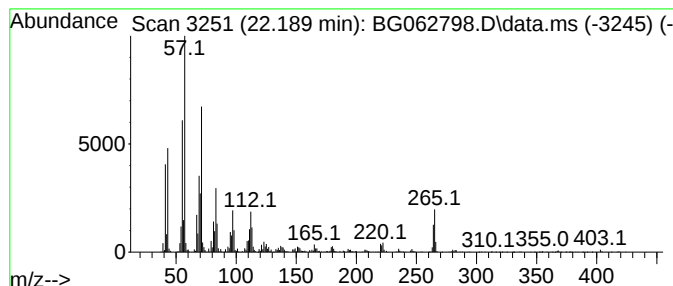
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 73 9-Octadecenoic acid, (E)- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.189	21.88 ng/ul	3605540	Chrysene-d12	21.531

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	53
2			2-Nonanol oleate	408	C27H52O2	1096360-61-0	49
3			Dimethylmalonic acid, 2,5-dichlo...	528	C29H46Cl2O4	1000363-69-5	38
4			Z-5-Nonadecene	266	C19H38	1000131-11-8	35
5			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062798.D
Acq On : 13 Sep 2024 21:36
Operator : JU/RC
Sample : P3898-20
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59

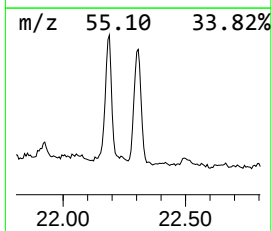
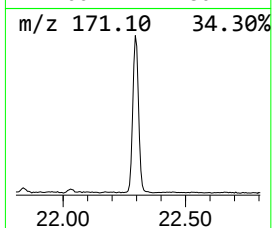
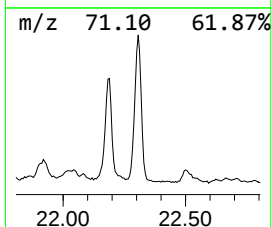
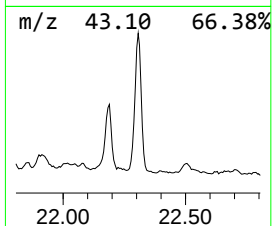
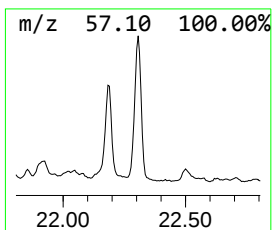
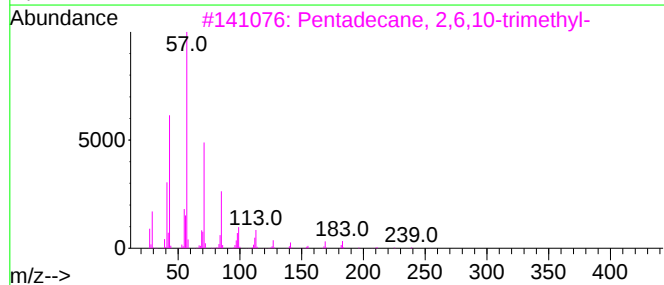
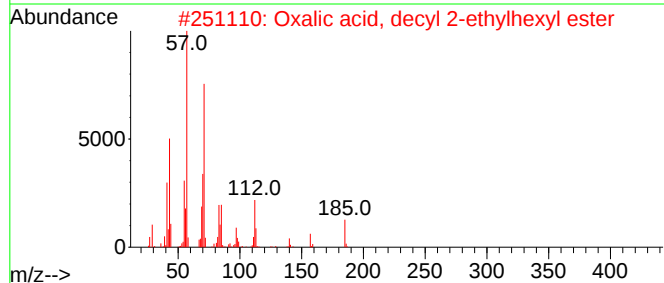
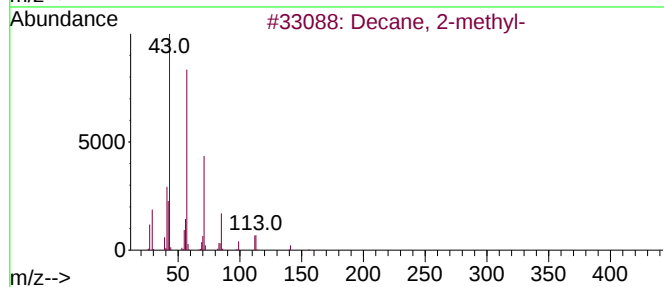
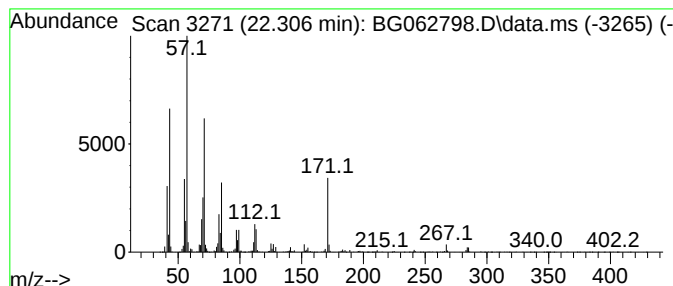
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.306	31.58 ng/ul	5205170	Chrysene-d12	21.531

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2-methyl-	156	C11H24	006975-98-0	70
2			Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	64
3			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64
4			Nonadecane	268	C19H40	000629-92-5	60
5			Pentacosane	352	C25H52	000629-99-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062798.D
 Acq On : 13 Sep 2024 21:36
 Operator : JU/RC
 Sample : P3898-20
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC59

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.937	28.4	ng/ul	10774200	1	7.917	7581300	20.0
Benzene, (1-met...	6.407	9.2	ng/ul	3490370	1	7.917	7581300	20.0
(DEL) Alkane: S...	6.478	9.7	ng/ul	3675990	1	7.917	7581300	20.0
(DEL) Alkane: C...	6.560	11.3	ng/ul	4273730	1	7.917	7581300	20.0
(DEL) Alkane: S...	6.918	24.1	ng/ul	9145270	1	7.917	7581300	20.0
(DEL) Alkane: S...	6.977	18.3	ng/ul	6931680	1	7.917	7581300	20.0
Benzene, 1,2,3-...	7.012	16.6	ng/ul	6278650	1	7.917	7581300	20.0
Benzene, 1-ethy...	7.306	14.3	ng/ul	5417710	1	7.917	7581300	20.0
2-Heptanone, 4,...	7.353	20.7	ng/ul	7852190	1	7.917	7581300	20.0
(DEL) Alkane: S...	7.576	97.5	ng/ul	36978600	1	7.917	7581300	20.0
(DEL) Alkane: S...	7.753	8.0	ng/ul	3047020	1	7.917	7581300	20.0
(DEL) Alkane: S...	7.847	8.9	ng/ul	3368250	1	7.917	7581300	20.0
1-Hexanol, 2-et...	8.011	53.0	ng/ul	20087600	1	7.917	7581300	20.0
(DEL) Alkane: S...	8.123	8.0	ng/ul	3014070	1	7.917	7581300	20.0
1,1-Hexylenedio...	8.152	14.6	ng/ul	5533710	1	7.917	7581300	20.0
Cyclohexanone, ...	8.364	20.8	ng/ul	7873070	1	7.917	7581300	20.0
(DEL) Alkane: S...	8.464	21.3	ng/ul	8061470	1	7.917	7581300	20.0
(DEL) Alkane: S...	8.522	9.4	ng/ul	3561060	1	7.917	7581300	20.0
Benzene, 1,3-di...	8.575	17.3	ng/ul	6561300	1	7.917	7581300	20.0
(DEL) Alkane: S...	8.699	21.0	ng/ul	7959450	1	7.917	7581300	20.0
Benzene, 1-meth...	8.728	13.4	ng/ul	5081490	1	7.917	7581300	20.0
Benzene, 1-ethy...	8.881	11.9	ng/ul	4490620	1	7.917	7581300	20.0
Benzene, 1-ethy...	8.928	17.7	ng/ul	6721870	1	7.917	7581300	20.0
Carbonic acid, ...	9.186	101.6	ng/ul	38528200	1	7.917	7581300	20.0
unknown-01	9.280	29.1	ng/ul	11014300	1	7.917	7581300	20.0
(DEL) Alkane: C...	9.856	3.6	ng/ul	8195450	2	10.714	45431500	20.0
(DEL) Alkane: S...	10.003	3.9	ng/ul	8734640	2	10.714	45431500	20.0
(DEL) Alkane: S...	10.150	3.7	ng/ul	8378390	2	10.714	45431500	20.0
(DEL) Alkane: S...	11.231	4.9	ng/ul	11120900	2	10.714	45431500	20.0
(DEL) Alkane: S...	11.754	5.1	ng/ul	11605500	2	10.714	45431500	20.0
(DEL) Alkane: S...	12.159	7.9	ng/ul	17867700	2	10.714	45431500	20.0
(DEL) Alkane: C...	12.794	16.2	ng/ul	4768820	3	14.551	5893380	20.0
1-Iodo-2-methyl...	12.947	15.4	ng/ul	4523300	3	14.551	5893380	20.0
unknown-02	13.305	13.1	ng/ul	3847550	3	14.551	5893380	20.0
Propanoic acid, ...	13.522	25.7	ng/ul	7578080	3	14.551	5893380	20.0
unknown-03	13.605	20.4	ng/ul	5999760	3	14.551	5893380	20.0
Naphthalene, 2, ...	13.734	20.7	ng/ul	6098290	3	14.551	5893380	20.0
Butanoic acid, ...	13.822	23.1	ng/ul	6804010	3	14.551	5893380	20.0
Naphthalene, 1, ...	13.881	26.0	ng/ul	7665260	3	14.551	5893380	20.0
(DEL) Alkane: S...	14.039	24.8	ng/ul	7312680	3	14.551	5893380	20.0
(DEL) Alkane: S...	14.075	15.7	ng/ul	4636180	3	14.551	5893380	20.0
unknown-04	14.180	31.8	ng/ul	9379490	3	14.551	5893380	20.0
(DEL) Alkane: S...	14.474	82.6	ng/ul	24327100	3	14.551	5893380	20.0
1H-Indole, 4-(3...	14.703	41.9	ng/ul	12355200	3	14.551	5893380	20.0
Naphthalene, 1, ...	15.032	12.7	ng/ul	3735100	3	14.551	5893380	20.0
(DEL) Alkane: S...	15.073	10.4	ng/ul	3064670	3	14.551	5893380	20.0
Naphthalene, 2, ...	15.326	49.8	ng/ul	14672000	3	14.551	5893380	20.0
(DEL) Alkane: S...	15.414	51.3	ng/ul	15114200	3	14.551	5893380	20.0
(DEL) Alkane: S...	15.808	13.3	ng/ul	3914160	3	14.551	5893380	20.0
(DEL) Alkane: S...	15.902	12.1	ng/ul	3560130	3	14.551	5893380	20.0
unknown-05	16.178	23.3	ng/ul	6385310	4	17.341	5489380	20.0
(DEL) Alkane: S...	16.272	48.4	ng/ul	13291000	4	17.341	5489380	20.0

unknown-06	16.607	10.9	ng/ul	2999750	4	17.341	5489380	20.0
2-Amino-5-isopr...	17.600	20.0	ng/ul	5489380	4	17.341	5489380	20.0
(DEL) Alkane: S...	17.800	34.6	ng/ul	9492230	4	17.341	5489380	20.0
Tridecanoic aci...	17.964	21.6	ng/ul	5927850	4	17.341	5489380	20.0
(DEL) Alkane: S...	18.487	23.4	ng/ul	6415770	4	17.341	5489380	20.0
(DEL) Alkane: S...	19.116	29.3	ng/ul	8038340	4	17.341	5489380	20.0
(DEL) Alkane: S...	20.215	23.3	ng/ul	3840190	5	21.531	3296200	20.0
(DEL) Alkane: S...	20.708	17.6	ng/ul	2901600	5	21.531	3296200	20.0
Oxalic acid, de...	21.202	44.4	ng/ul	7313450	5	21.531	3296200	20.0
(DEL) Alkane: S...	21.724	20.0	ng/ul	3296200	5	21.531	3296200	20.0
9-Octadecenoic ...	22.189	21.9	ng/ul	3605540	5	21.531	3296200	20.0
(DEL) Alkane: S...	22.306	31.6	ng/ul	5205170	5	21.531	3296200	20.0

Instrument :

BNA_G

ClientSampleId :

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