

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
 Data File : BG062769.D
 Acq On : 12 Sep 2024 19:42
 Operator : JU/RC
 Sample : P3877-11DL2 30X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCVY4DL2

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: BG062769.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.931	477	484	491	rVB2	139793	217670	5.48%	1.021%
2	6.189	522	528	535	rBV4	21333	51700	1.30%	0.243%
3	6.466	570	575	580	rBV	40059	58339	1.47%	0.274%
4	6.548	580	589	593	rBV3	33403	79412	2.00%	0.373%
5	6.900	645	649	654	rVB2	55026	89210	2.25%	0.419%
6	6.959	654	659	662	rBV	64256	90897	2.29%	0.427%
7	6.994	662	665	670	rVB	53821	78578	1.98%	0.369%
8	7.065	670	677	683	rVB4	81654	160392	4.04%	0.753%
9	7.130	683	688	693	rVB	40805	57707	1.45%	0.271%
10	7.294	710	716	719	rBV2	36656	61067	1.54%	0.287%
11	7.417	730	737	743	rVB4	28160	59178	1.49%	0.278%
12	7.535	751	757	766	rBV2	326150	602217	15.16%	2.826%
13	7.823	798	806	811	rBV7	17950	48024	1.21%	0.225%
14	7.899	811	819	825	rBV2	168709	338926	8.53%	1.590%
15	7.964	825	830	836	rBV	78273	127088	3.20%	0.596%
16	8.023	836	840	848	rVB2	32623	61471	1.55%	0.288%
17	8.193	866	869	876	rVB4	31918	55337	1.39%	0.260%
18	8.440	906	911	917	rVB3	44879	99445	2.50%	0.467%
19	8.499	917	921	924	rVB	36885	49585	1.25%	0.233%
20	8.563	924	932	936	rBV3	66603	171785	4.32%	0.806%
21	8.669	945	950	954	rBV	57478	94572	2.38%	0.444%
22	8.710	954	957	965	rVB3	33791	63370	1.60%	0.297%
23	8.904	985	990	997	rVB	34655	63080	1.59%	0.296%
24	9.004	997	1007	1018	rBV3	48392	131856	3.32%	0.619%
25	9.133	1018	1029	1035	rVV	316094	502328	12.64%	2.357%
26	9.256	1039	1050	1057	rBV9	38162	111623	2.81%	0.524%
27	9.691	1116	1124	1131	rVB3	26339	53889	1.36%	0.253%
28	9.785	1131	1140	1143	rBV7	22179	46902	1.18%	0.220%
29	9.838	1143	1149	1155	rBV6	38831	85385	2.15%	0.401%
30	9.979	1165	1173	1178	rVB	40792	83459	2.10%	0.392%
31	10.050	1178	1185	1189	rVV4	26947	45539	1.15%	0.214%
32	10.126	1195	1198	1202	rVV2	35548	60870	1.53%	0.286%
33	10.232	1209	1216	1220	rBV2	24716	46010	1.16%	0.216%
34	10.390	1238	1243	1248	rBV2	30761	53310	1.34%	0.250%
35	10.690	1282	1294	1300	rBV2	281733	732835	18.45%	3.439%
36	10.743	1300	1303	1310	rVB2	33874	57611	1.45%	0.270%

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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
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37	10.813	1310	1315	1320	rBV3	59880	106856	2.69%	0.501%
38	10.860	1320	1323	1329	rVB	29306	46300	1.17%	0.217%
39	11.072	1353	1359	1372	rBV	136901	225821	5.68%	1.060%
40	11.201	1372	1381	1387	rBV5	25314	61839	1.56%	0.290%

41	11.613	1442	1451	1456	rVV7	21753	52888	1.33%	0.248%
42	11.724	1464	1470	1474	rBV3	59955	97739	2.46%	0.459%
43	11.795	1474	1482	1489	rVB2	98416	194944	4.91%	0.915%
44	11.959	1502	1510	1516	rBV	77853	128672	3.24%	0.604%
45	12.118	1531	1537	1540	rBV	119576	190334	4.79%	0.893%

46	12.153	1540	1543	1548	rVB2	78572	109823	2.76%	0.515%
47	12.359	1572	1578	1587	rVB2	154691	271554	6.84%	1.274%
48	12.570	1610	1614	1623	rVB3	35425	74210	1.87%	0.348%
49	12.758	1637	1646	1650	rBV3	28854	66376	1.67%	0.311%
50	12.929	1670	1675	1681	rBV4	26949	46609	1.17%	0.219%

51	13.076	1695	1700	1707	rVV2	39377	64001	1.61%	0.300%
52	13.358	1742	1748	1756	rVV	151618	223346	5.62%	1.048%
53	13.487	1767	1770	1776	rVV5	20447	45566	1.15%	0.214%
54	13.581	1780	1786	1793	rVB9	20086	52055	1.31%	0.244%
55	13.710	1799	1808	1814	rBV6	43679	110389	2.78%	0.518%

56	13.792	1817	1822	1827	rBV2	35051	51629	1.30%	0.242%
57	13.857	1827	1833	1835	rBV3	39739	80207	2.02%	0.376%
58	14.021	1853	1861	1865	rVB2	58521	99709	2.51%	0.468%
59	14.163	1881	1885	1893	rVB2	66776	86672	2.18%	0.407%
60	14.257	1898	1901	1911	rVB8	21605	57415	1.45%	0.269%

61	14.433	1926	1931	1936	rVB	349749	450141	11.33%	2.112%
62	14.533	1940	1948	1954	rVB	552354	886569	22.32%	4.160%
63	14.603	1954	1960	1967	rBV8	48899	92287	2.32%	0.433%
64	14.674	1967	1972	1979	rVB	210124	301158	7.58%	1.413%
65	14.738	1979	1983	1991	rVB7	44470	95064	2.39%	0.446%

66	14.926	2013	2015	2020	rVB4	38507	48662	1.22%	0.228%
67	15.056	2033	2037	2041	rBV4	31506	52071	1.31%	0.244%
68	15.091	2041	2043	2047	rVV4	36766	59327	1.49%	0.278%
69	15.155	2049	2054	2061	rVV	179871	338012	8.51%	1.586%
70	15.220	2062	2065	2069	rVB2	52415	67765	1.71%	0.318%

71	15.291	2072	2077	2085	rVV	220163	326189	8.21%	1.531%
72	15.385	2085	2093	2098	rVB	161364	259492	6.53%	1.218%
73	15.625	2129	2134	2138	rBV	49773	77322	1.95%	0.363%
74	15.790	2157	2162	2170	rVB2	67019	113504	2.86%	0.533%
75	15.890	2174	2179	2185	rVB7	47408	80115	2.02%	0.376%

76	16.160	2220	2225	2230	rVV	91730	137254	3.45%	0.644%
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77	16.248	2235	2240	2242	rBV	168171	223228	5.62%	1.047%
78	16.460	2273	2276	2280	rVB3	38983	47231	1.19%	0.222%
79	16.589	2292	2298	2303	rBV3	61000	101692	2.56%	0.477%
80	16.930	2349	2356	2367	rBV	2620418	3972737	100.00%	18.642%
81	17.036	2371	2374	2379	rVV	155200	212506	5.35%	0.997%
82	17.094	2379	2384	2394	rVB3	90870	178769	4.50%	0.839%
83	17.276	2409	2415	2422	rBV	850100	1278439	32.18%	5.999%
84	17.371	2427	2431	2435	rVV	75955	115554	2.91%	0.542%
85	17.418	2435	2439	2442	rVB4	54960	69166	1.74%	0.325%
86	17.564	2459	2464	2471	rVB4	69938	133084	3.35%	0.624%
87	17.776	2496	2500	2505	rVB	143806	176855	4.45%	0.830%
88	17.946	2525	2529	2534	rVB2	44561	58145	1.46%	0.273%
89	18.469	2613	2618	2623	rBV	98443	129465	3.26%	0.608%
90	19.098	2721	2725	2727	rBV	64327	83479	2.10%	0.392%
91	19.210	2740	2744	2748	rBV3	39221	52077	1.31%	0.244%
92	19.674	2817	2823	2829	rBV2	72342	132311	3.33%	0.621%
93	20.208	2910	2914	2921	rBV4	48708	103185	2.60%	0.484%
94	20.702	2994	2998	3002	rBV2	47264	56493	1.42%	0.265%
95	21.195	3078	3082	3086	rVB	76223	93255	2.35%	0.438%
96	21.519	3131	3137	3144	rBV	972525	1511869	38.06%	7.094%
97	21.718	3165	3171	3176	rBV3	38446	63542	1.60%	0.298%
98	22.294	3264	3269	3276	rBV3	43380	77395	1.95%	0.363%
99	24.362	3618	3621	3630	rVB5	24738	59651	1.50%	0.280%
100	24.586	3649	3659	3673	rVB	602333	1660227	41.79%	7.791%

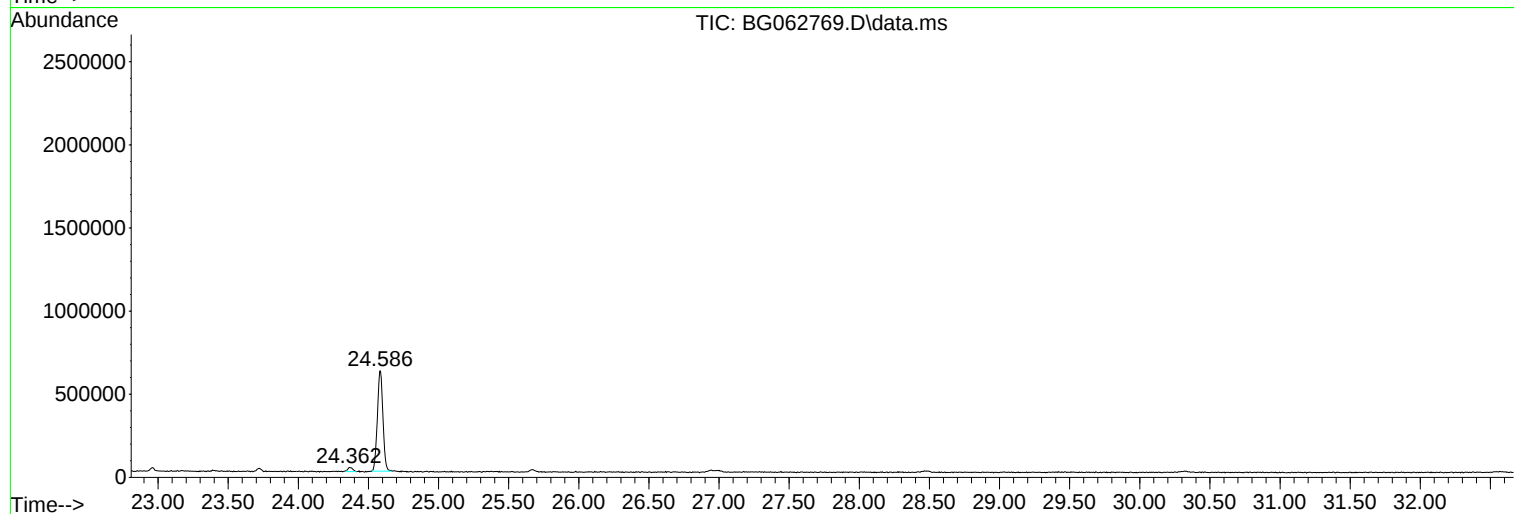
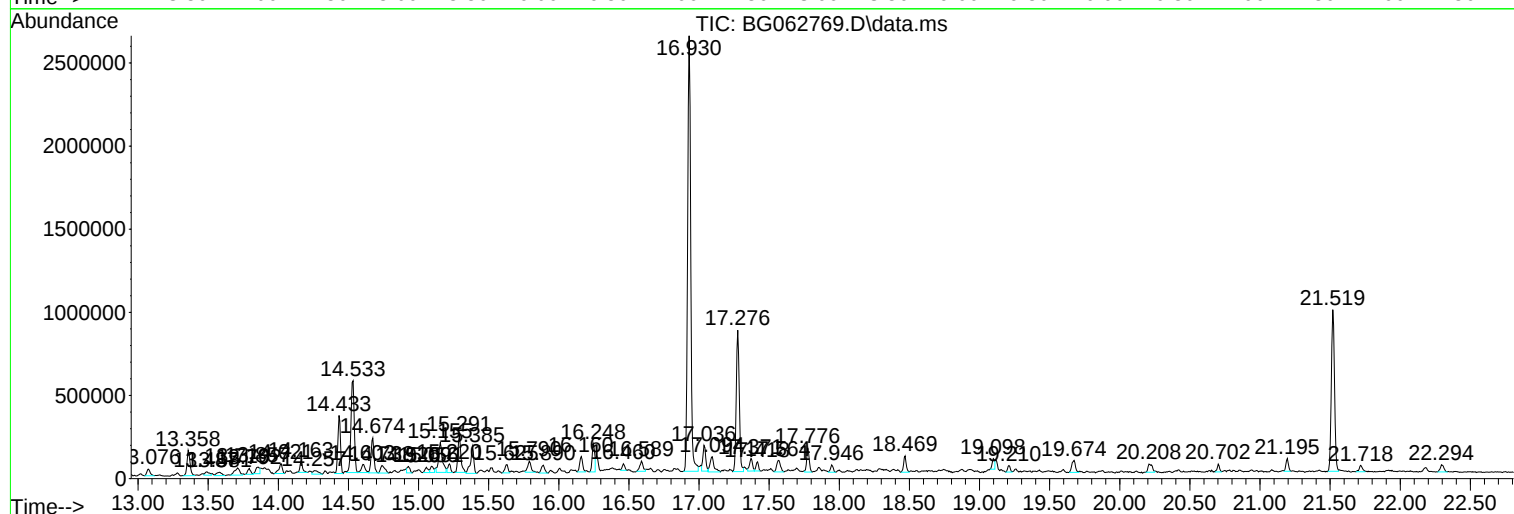
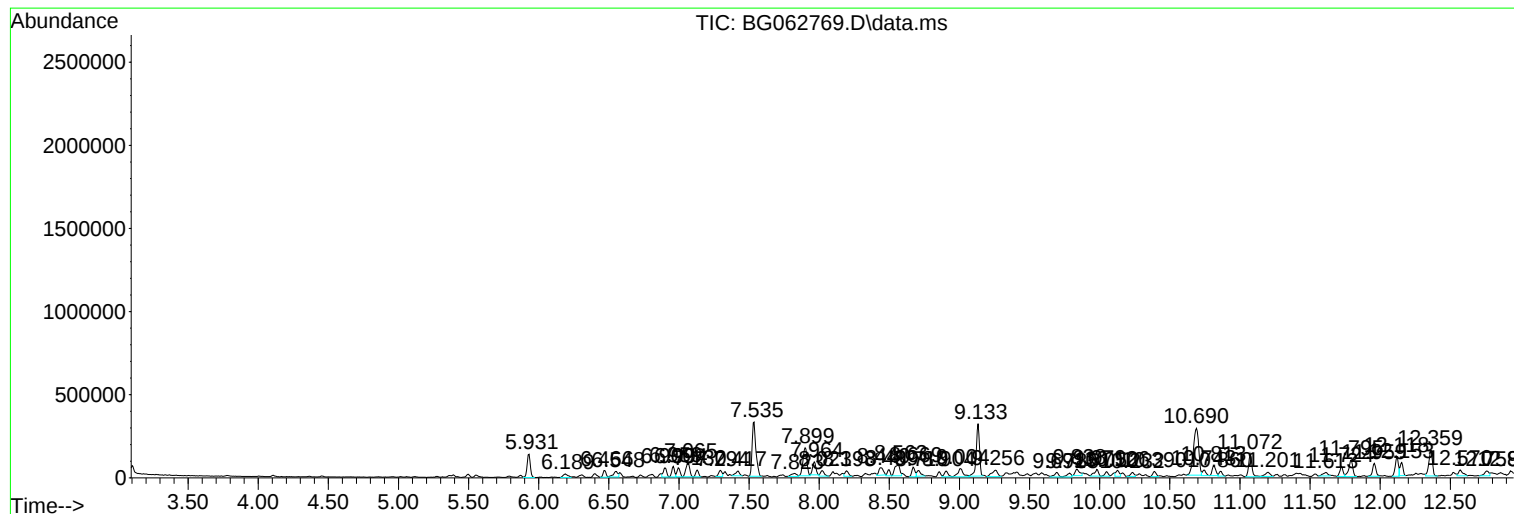
Sum of corrected areas: 21310908

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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



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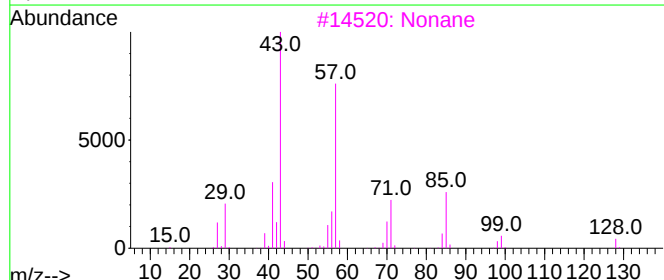
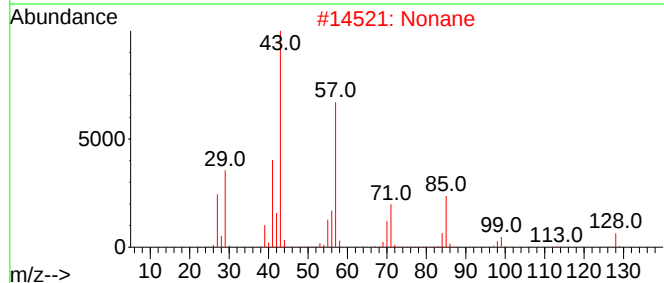
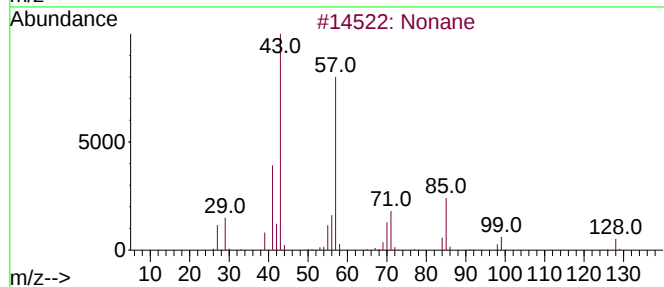
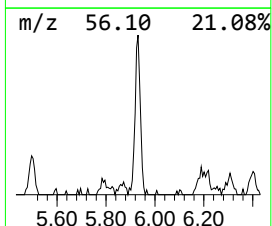
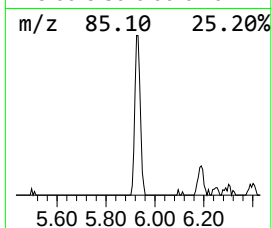
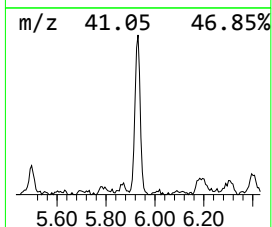
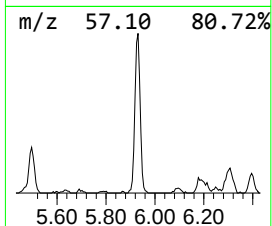
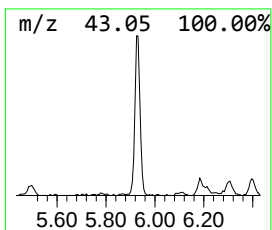
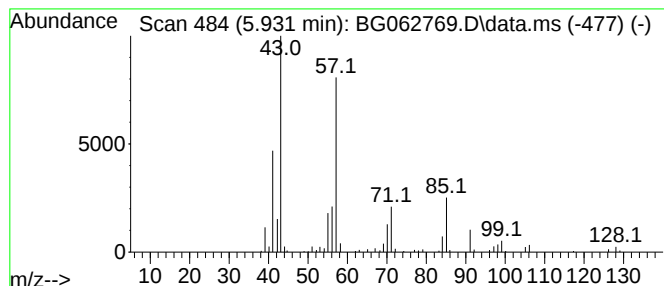
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.931	12.84 ng/ul	217670	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	90
2	Nonane			128	C9H20	000111-84-2	86
3	Nonane			128	C9H20	000111-84-2	78
4	Nonane			128	C9H20	000111-84-2	78
5	Decane			142	C10H22	000124-18-5	59



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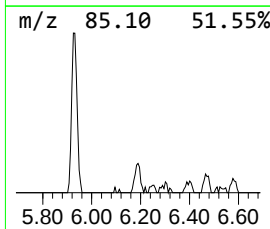
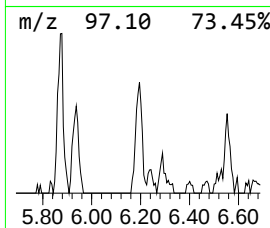
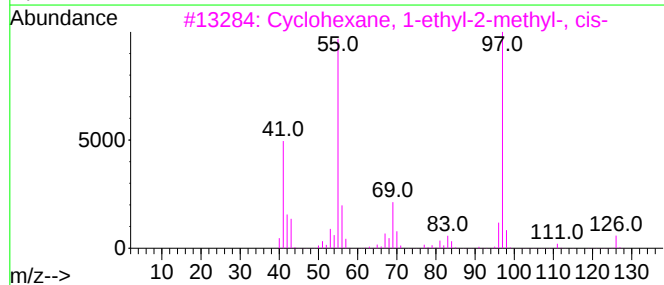
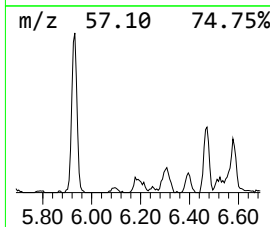
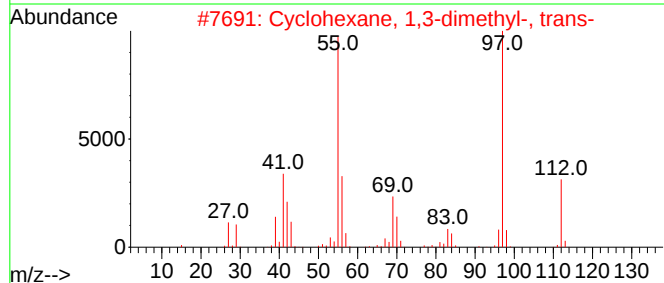
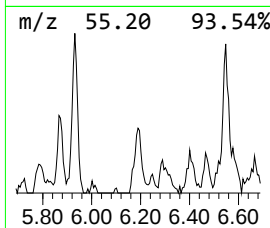
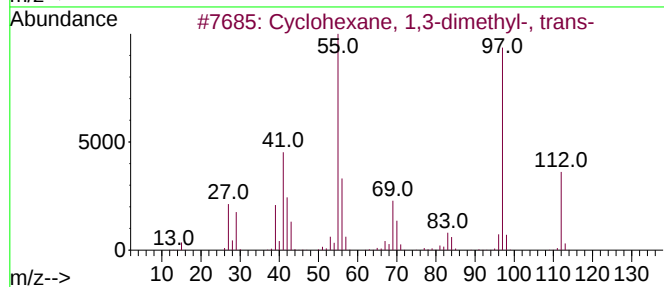
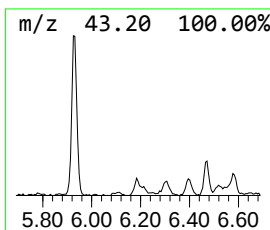
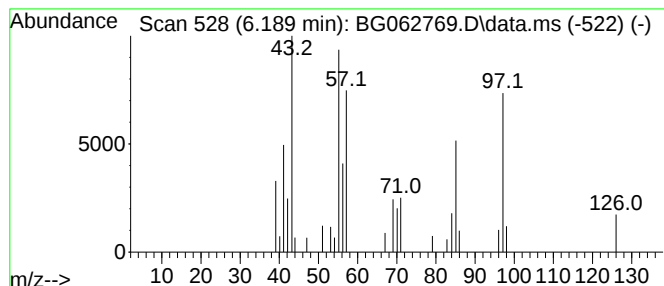
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Cyclic6.189 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.189	3.05 ng/ul	51700	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	47
2			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	47
3			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	43
4			1-Nonene	126	C9H18	000124-11-8	38
5			3-Heptene, 3-ethyl-	126	C9H18	074764-46-8	38



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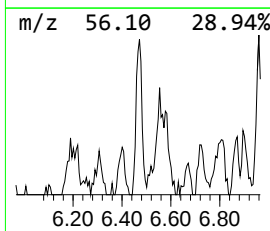
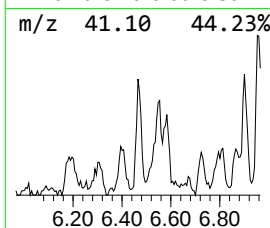
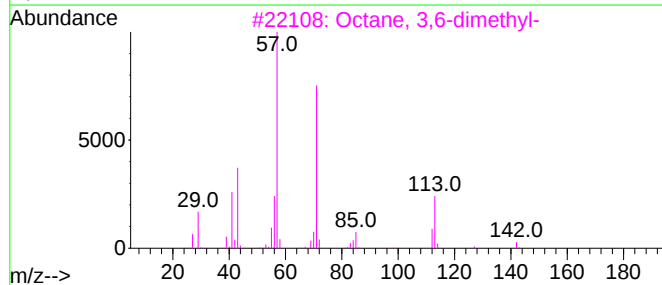
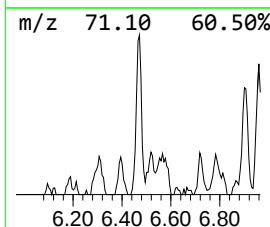
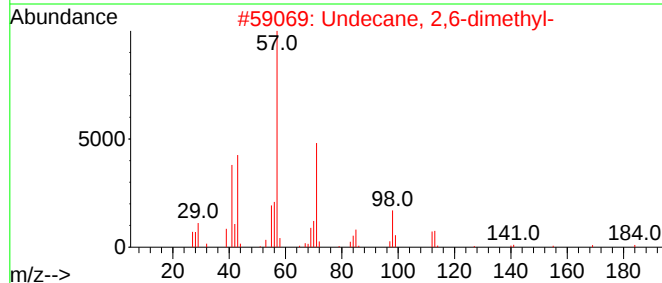
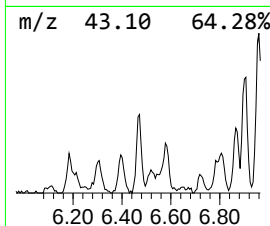
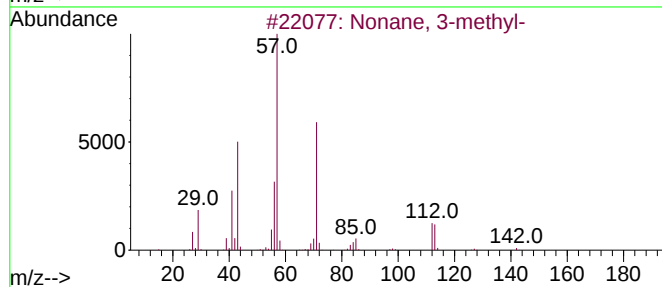
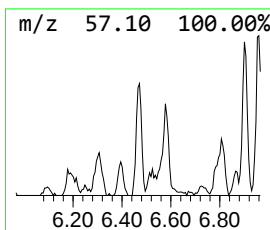
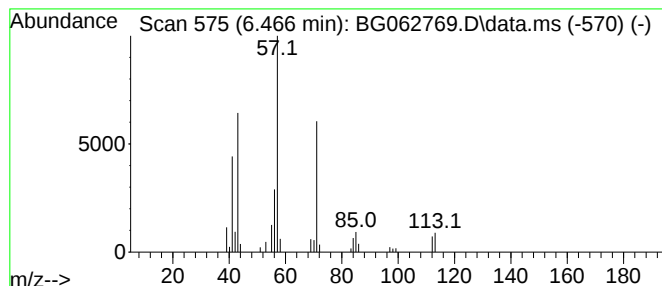
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ClientSampleId :
DCVY4DL2

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 3 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.466	3.44 ng/ul	58339	1,4-Dichlorobenzene-d4			7.905
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-		142	C10H22	005911-04-6	78
2	Undecane, 2,6-dimethyl-		184	C13H28	017301-23-4	72
3	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	72
4	Decane, 2,4-dimethyl-		170	C12H26	002801-84-5	72
5	Nonane, 3-methyl-		142	C10H22	005911-04-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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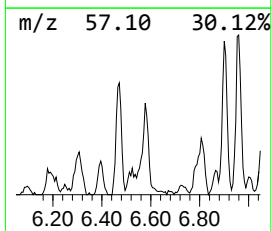
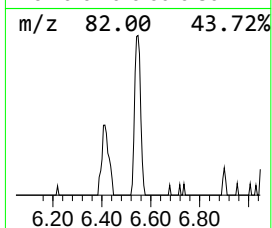
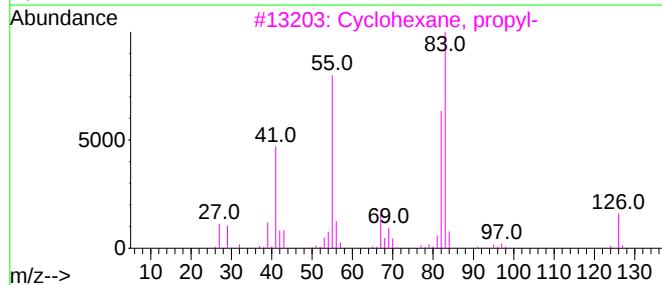
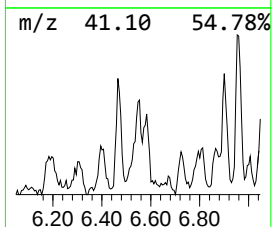
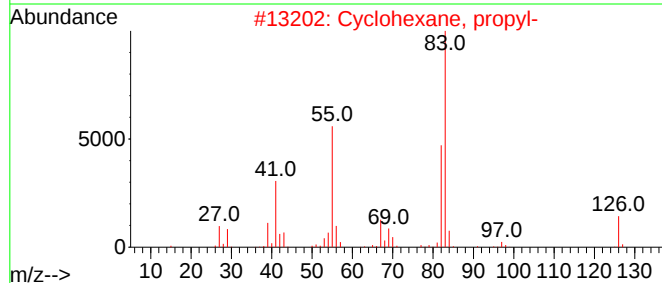
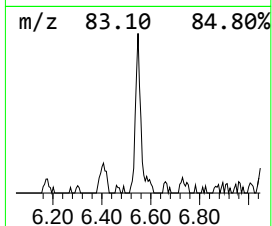
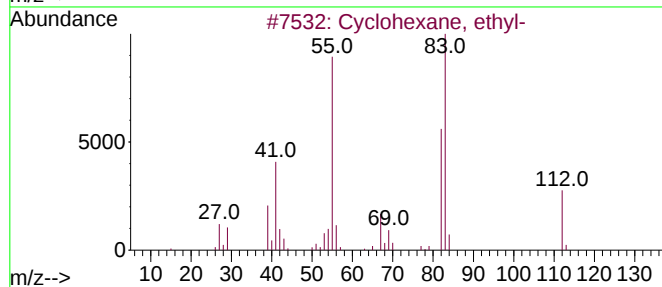
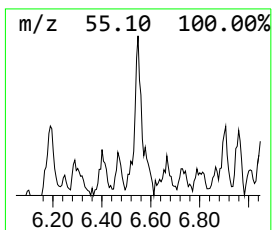
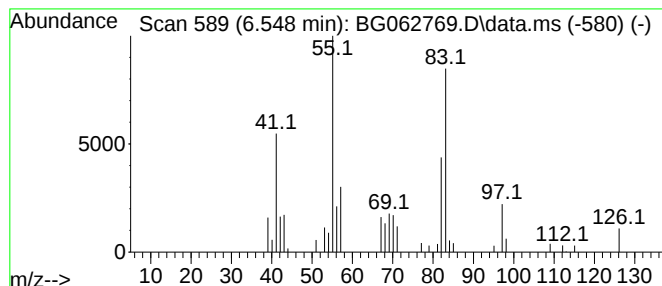
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic6.548 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.548	4.69 ng/ul	79412	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	58
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
5			Oxalic acid, cyclohexyl nonyl ester	298	C17H30O4	1000309-31-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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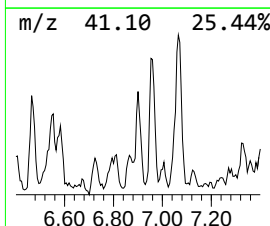
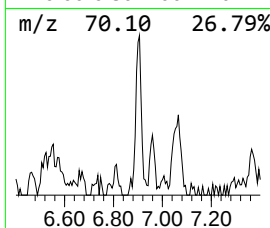
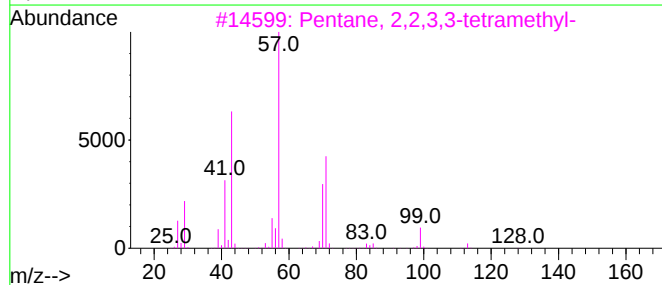
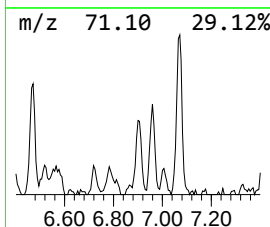
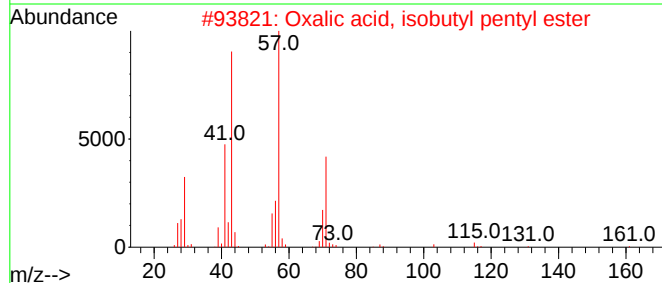
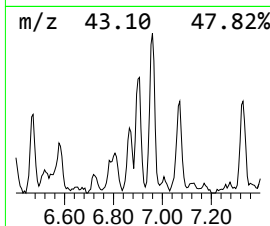
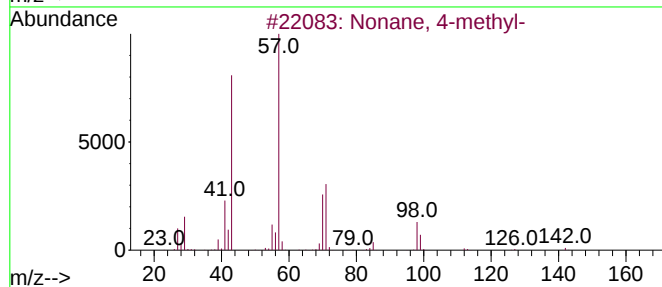
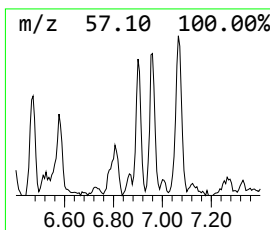
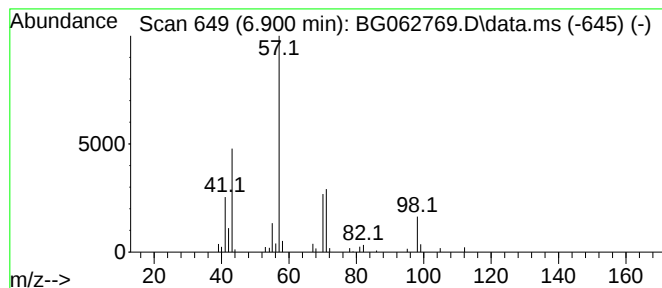
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.900	5.26 ng/ul	89210	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	50
2			Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	43
3			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	40
4			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	40
5			Dodecane, 6-methyl-	184	C13H28	006044-71-9	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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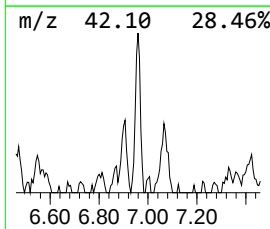
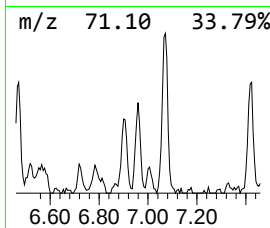
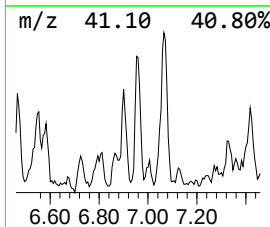
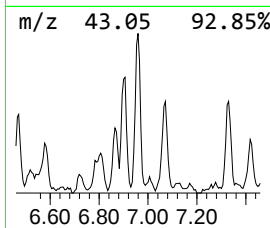
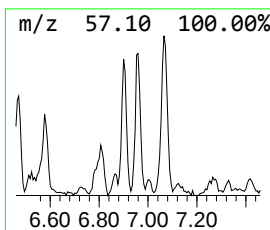
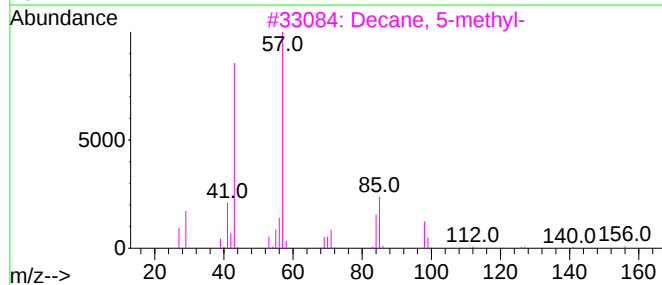
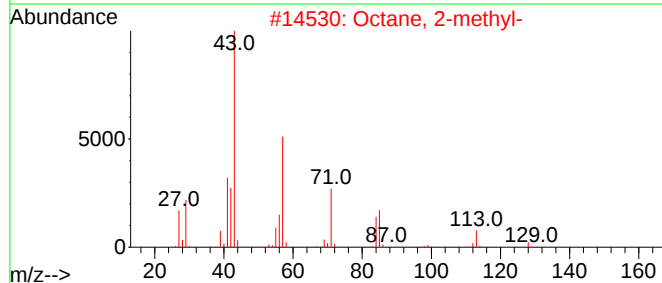
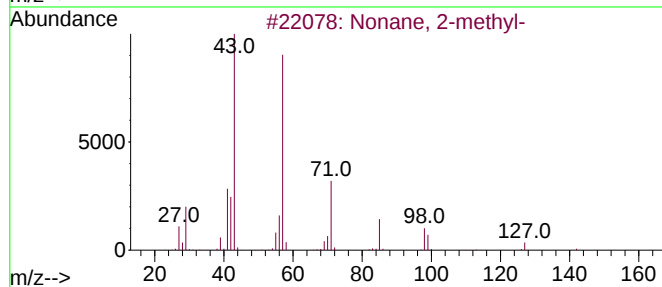
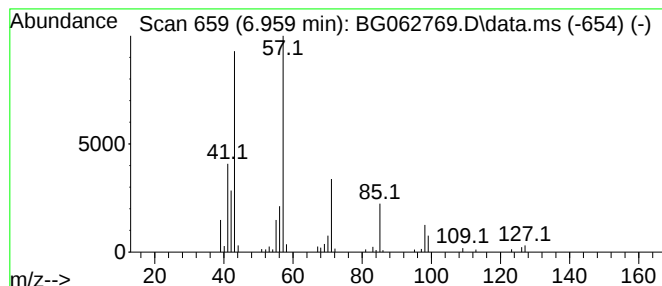
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.959	5.36 ng/ul	90897	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	64
2			Octane, 2-methyl-	128	C9H20	003221-61-2	53
3			Decane, 5-methyl-	156	C11H24	013151-35-4	53
4			Decane	142	C10H22	000124-18-5	50
5			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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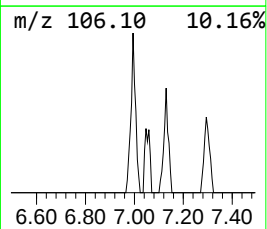
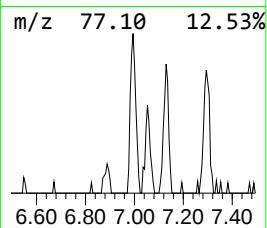
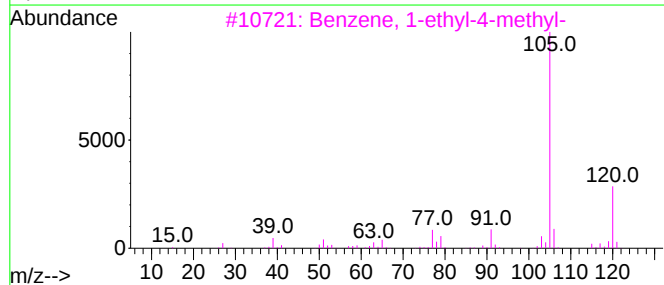
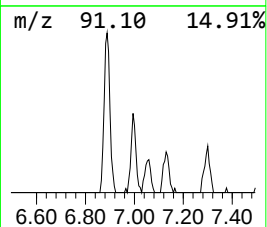
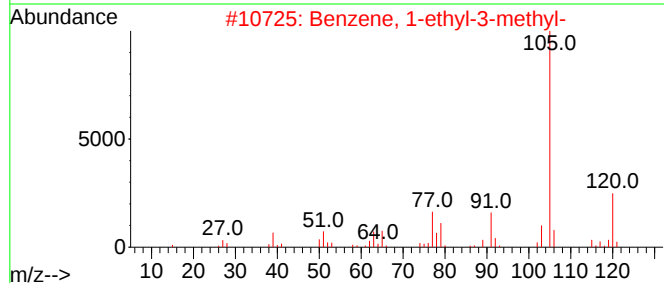
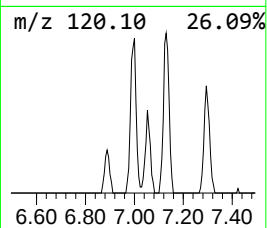
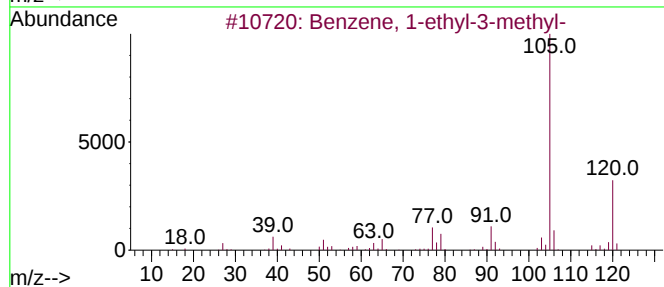
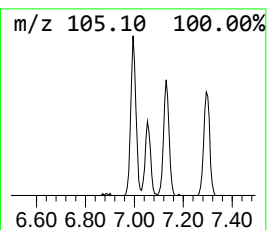
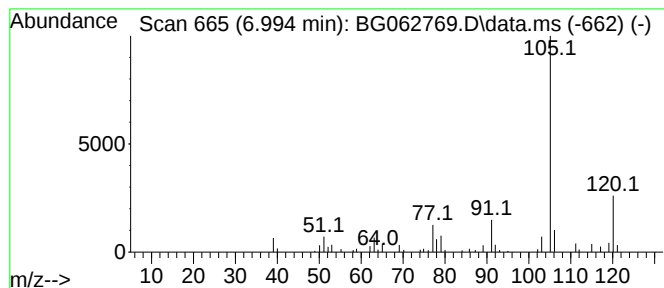
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.994	4.64 ng/ul	78578	1,4-Dichlorobenzene-d4	7.905

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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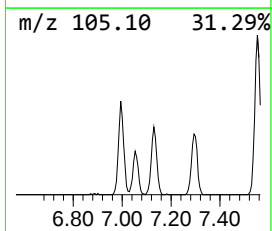
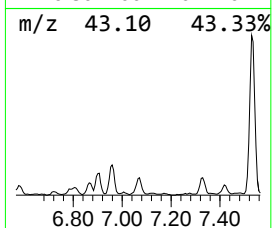
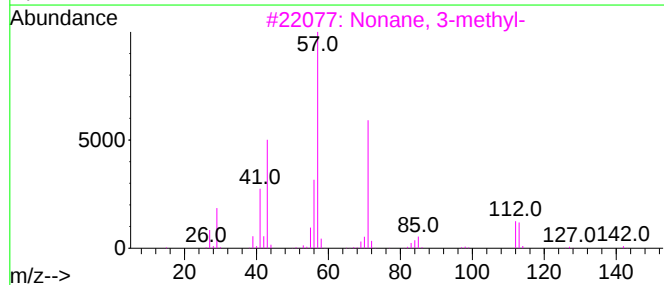
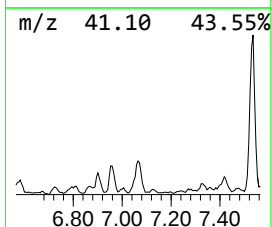
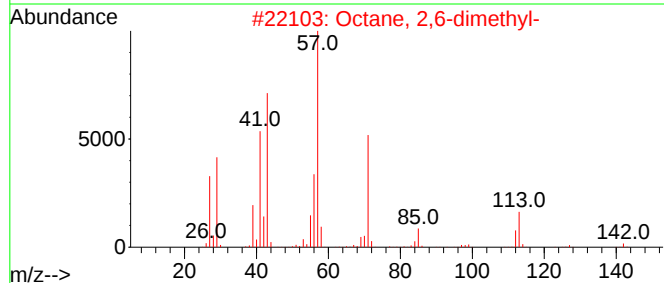
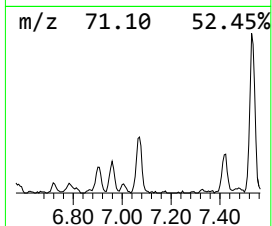
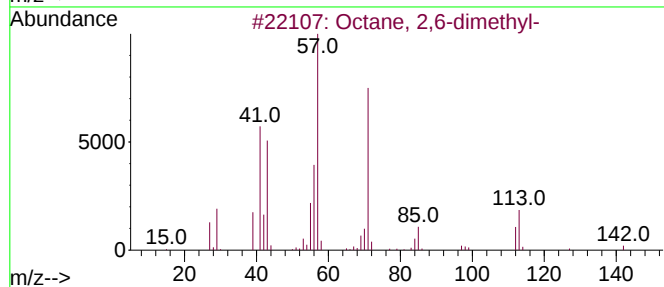
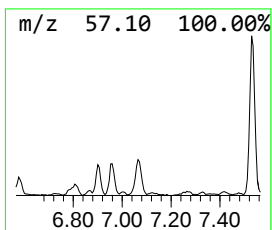
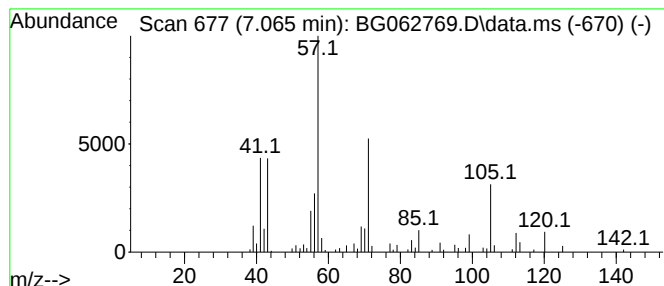
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.065	9.46 ng/ul	160392	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	58	
2	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	58	
3	Nonane, 3-methyl-		142	C10H22	005911-04-6	52	
4	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	52	
5	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	52	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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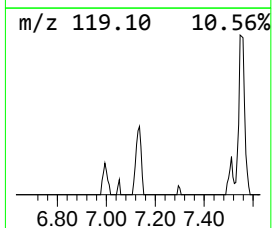
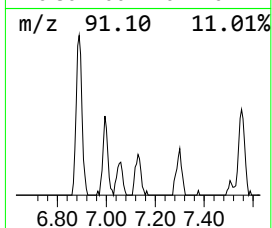
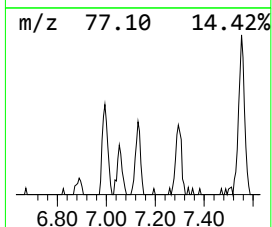
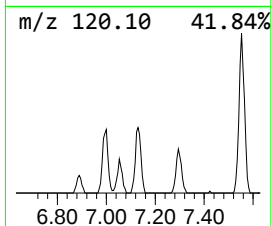
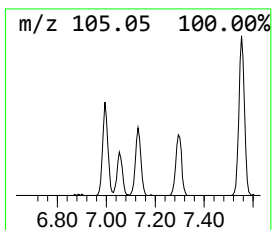
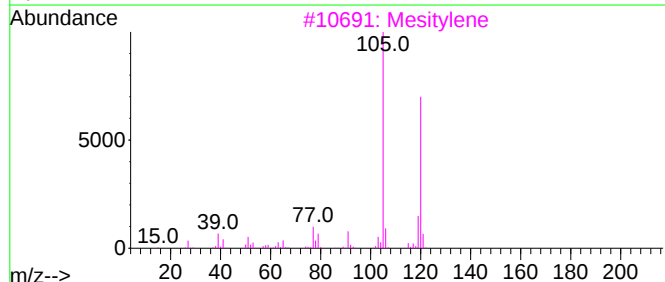
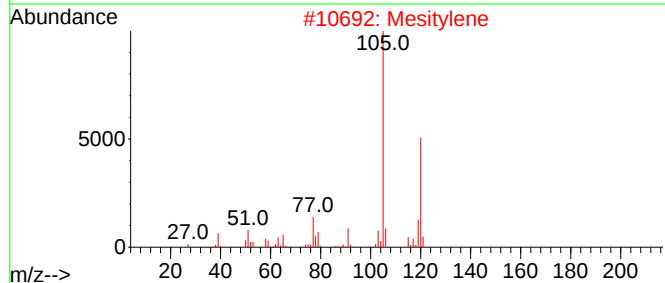
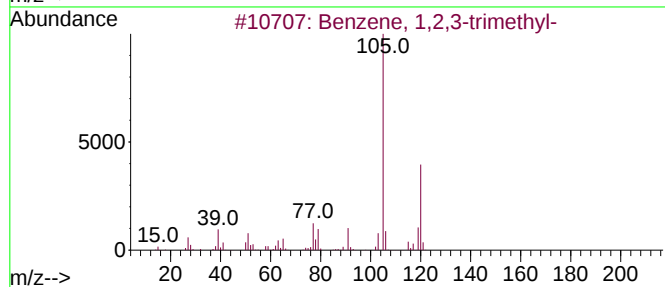
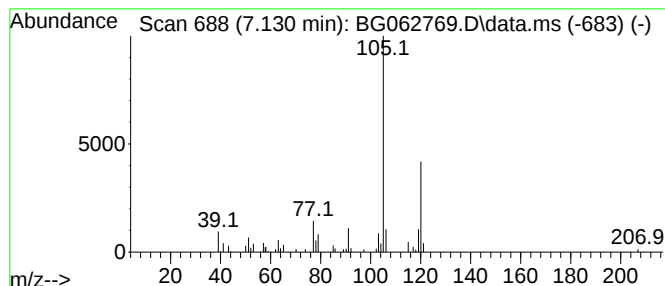
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.130	3.41 ng/ul	57707	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2			Mesitylene	120	C9H12	000108-67-8	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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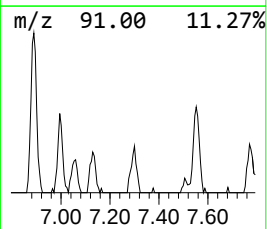
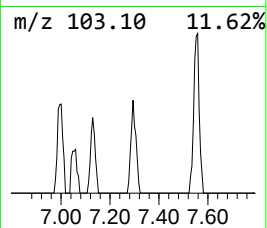
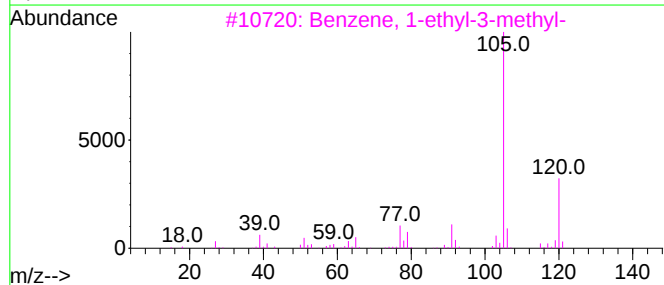
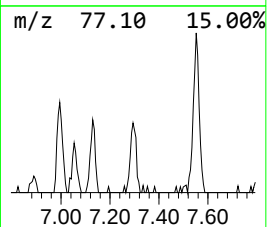
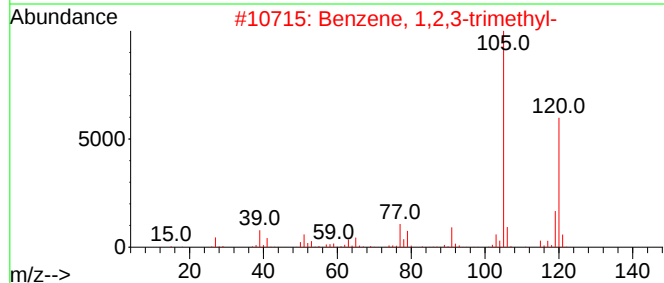
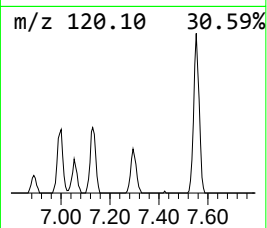
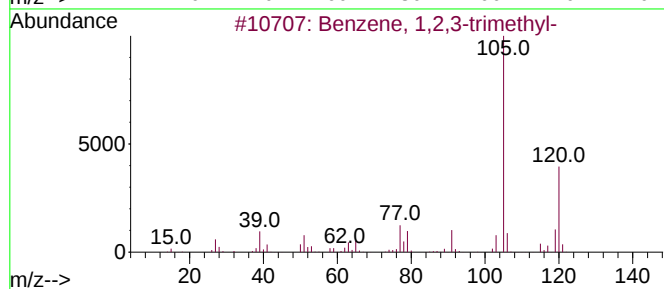
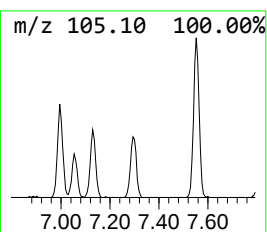
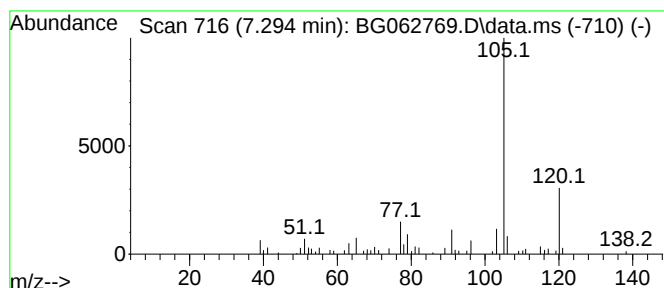
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2,4-Nonadiyne Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.294	3.60 ng/ul	61067	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
4			2,4-Nonadiyne	120	C9H12	063621-15-8	87
5			Mesitylene	120	C9H12	000108-67-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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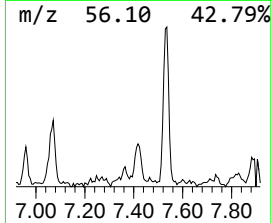
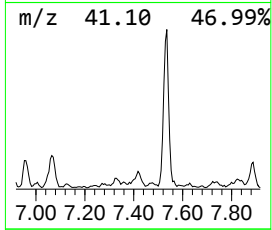
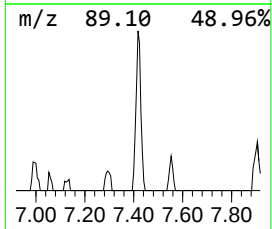
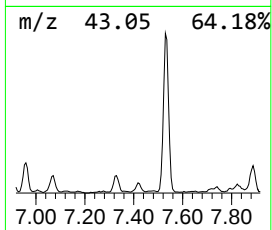
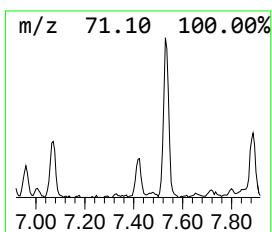
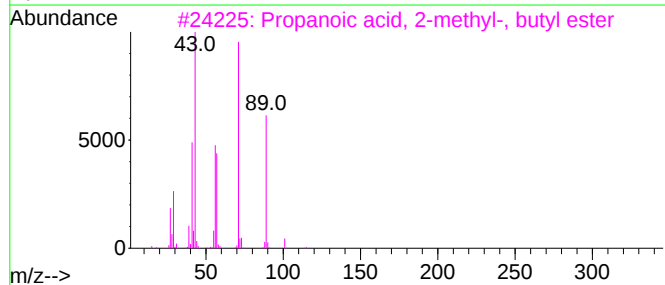
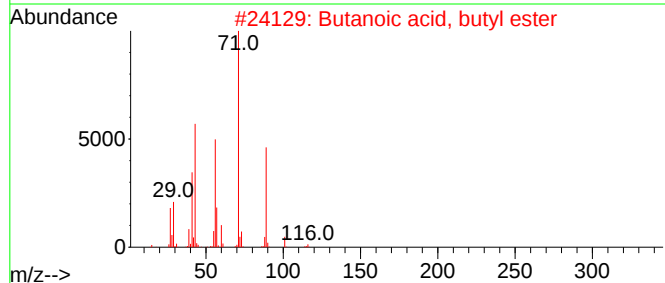
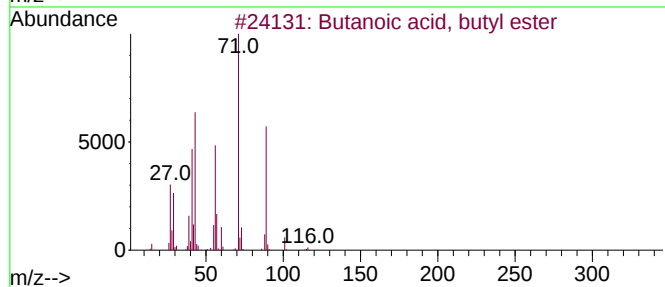
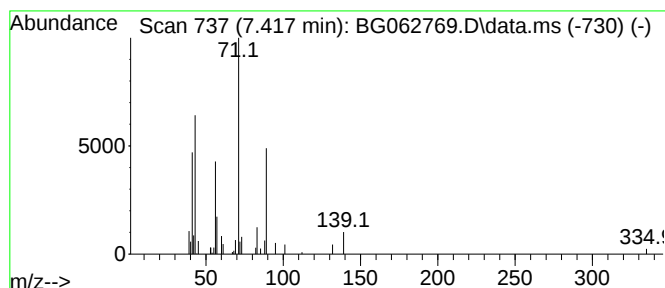
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Butanoic acid, butyl ester Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.417	3.49 ng/ul	59178	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	72
2			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	72
3			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	64
4			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	64
5			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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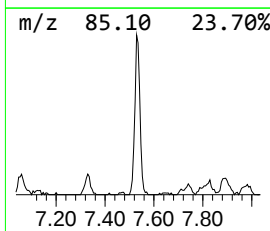
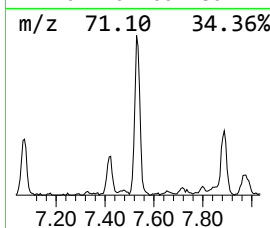
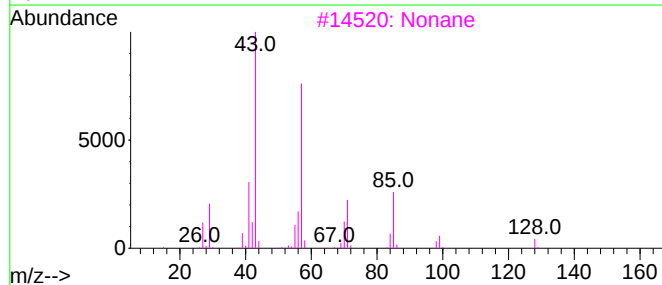
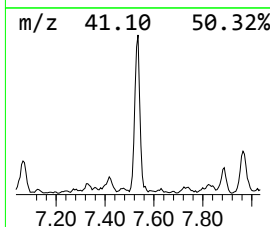
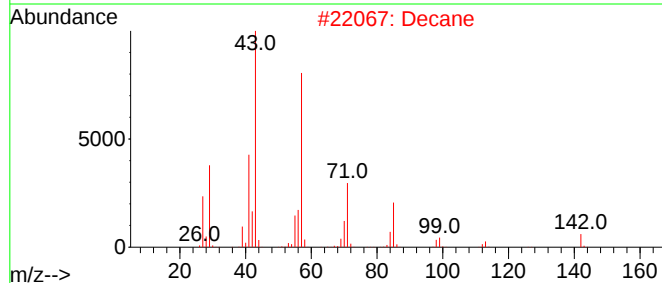
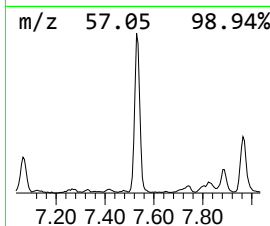
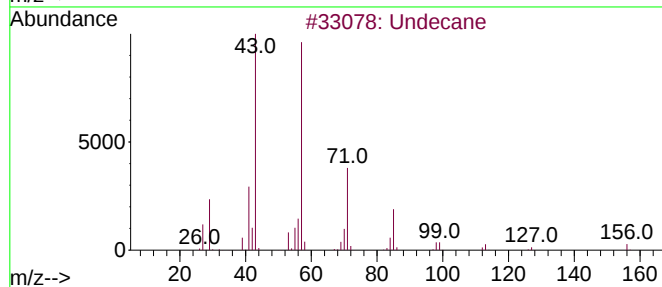
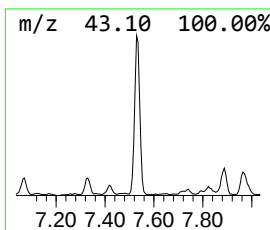
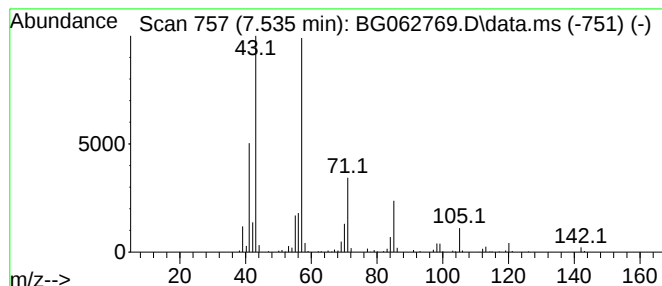
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.535	35.54 ng/ul	602217	1,4-Dichlorobenzene-d4	7.905

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane	156	C11H24	001120-21-4	80
2		Decane	142	C10H22	000124-18-5	80
3		Nonane	128	C9H20	000111-84-2	72
4		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
5		Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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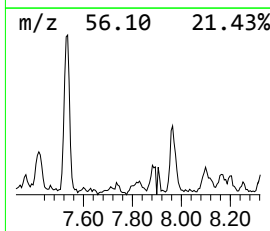
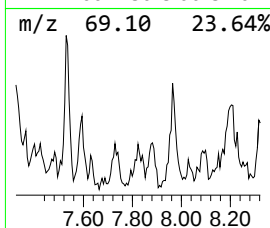
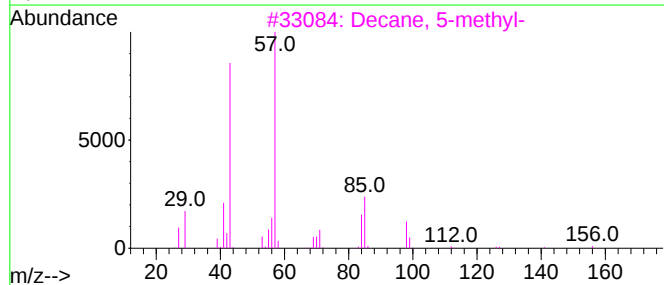
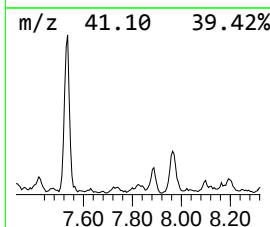
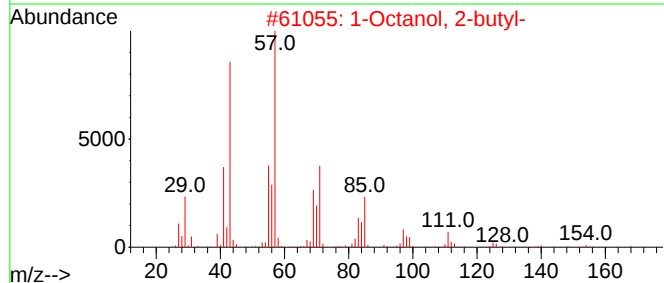
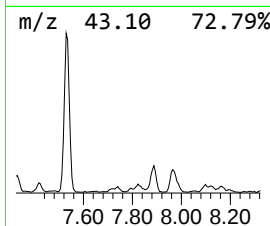
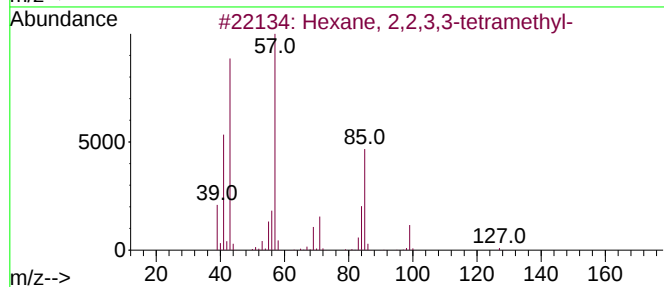
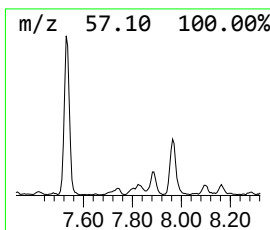
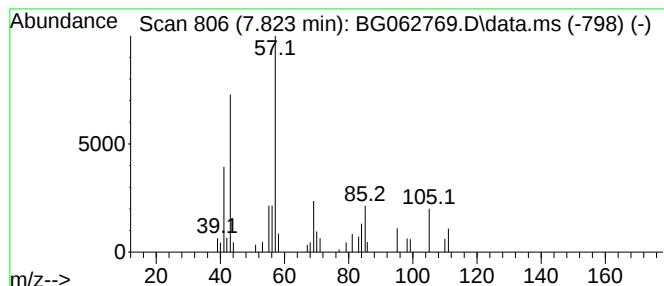
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.823	2.83 ng/ul	48024	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	43
2			1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	37
3			Decane, 5-methyl-	156	C11H24	013151-35-4	32
4			Decane	142	C10H22	000124-18-5	28
5			Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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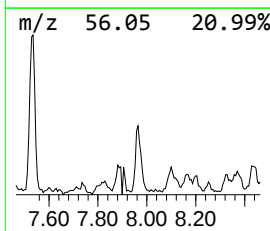
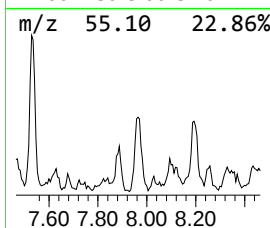
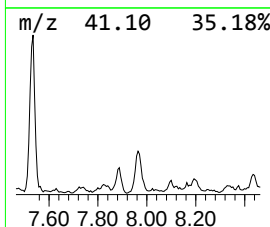
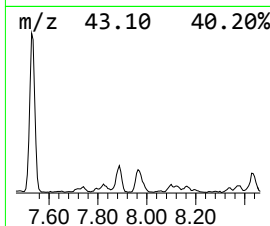
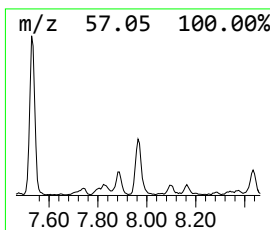
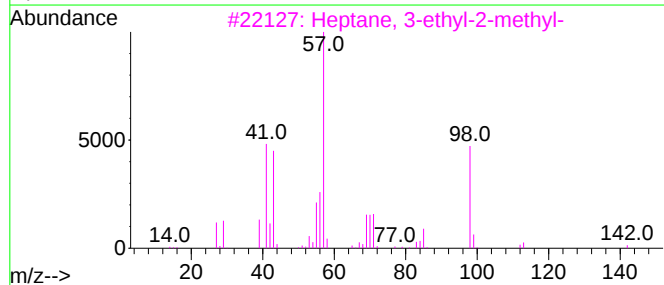
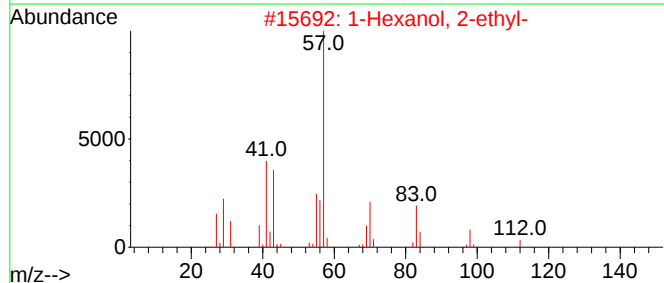
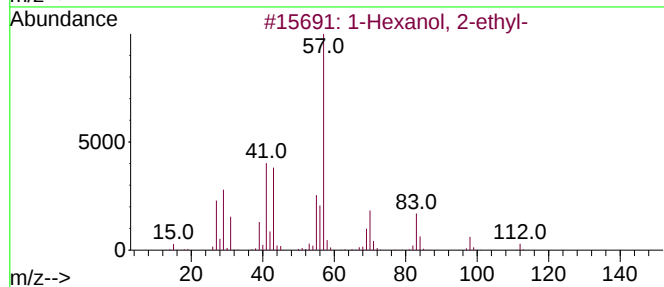
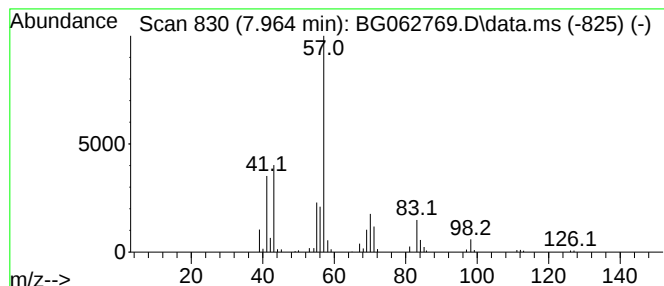
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1-Hexanol, 2-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.964	7.50 ng/ul	127088	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	Heptane, 3-ethyl-2-methyl-			142	C10H22	014676-29-0	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	59
5	1-Pentanol, 2-ethyl-4-methyl-			130	C8H18O	000106-67-2	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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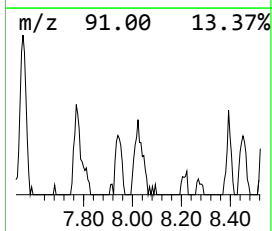
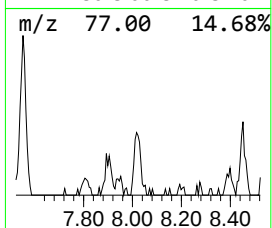
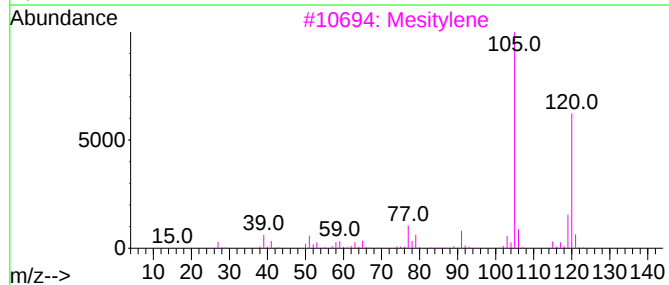
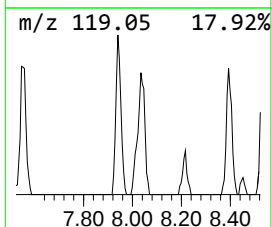
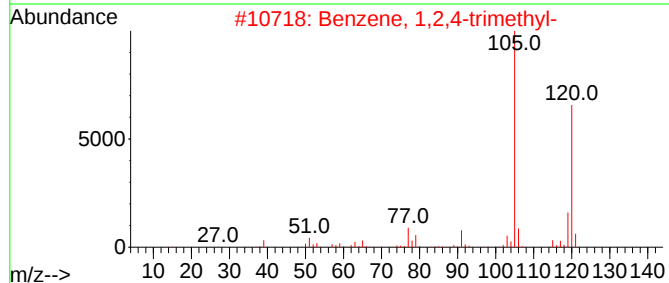
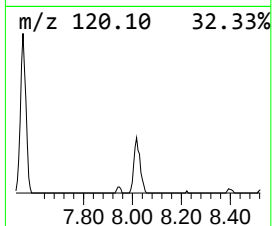
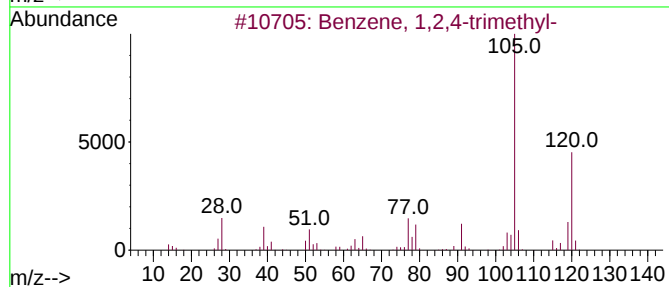
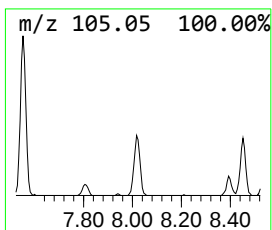
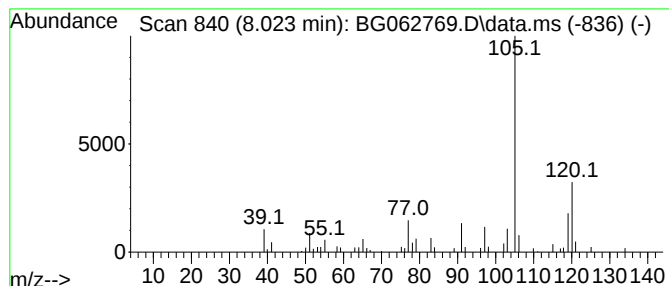
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,2,4-trimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.023	3.63 ng/ul	61471	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	80
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	80
3			Mesitylene	120	C9H12	000108-67-8	80
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
5			Mesitylene	120	C9H12	000108-67-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
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Misc :
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ClientSampleId :
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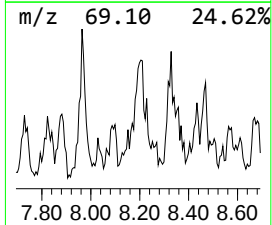
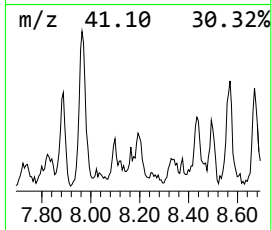
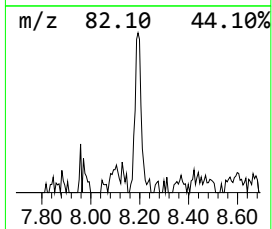
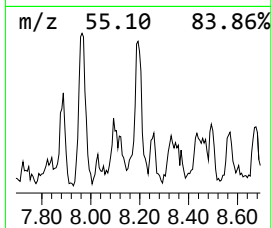
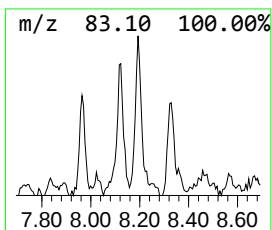
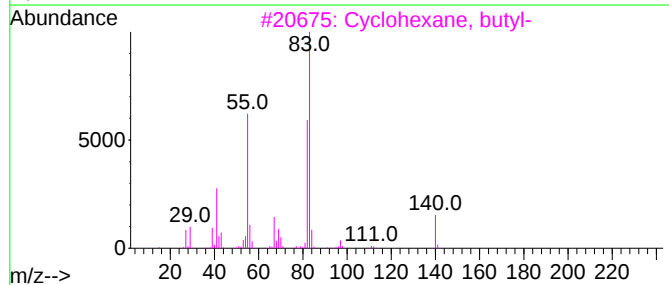
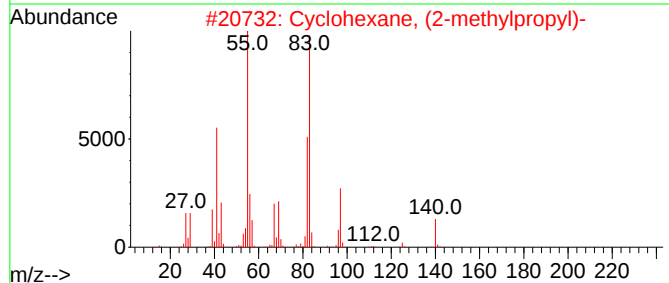
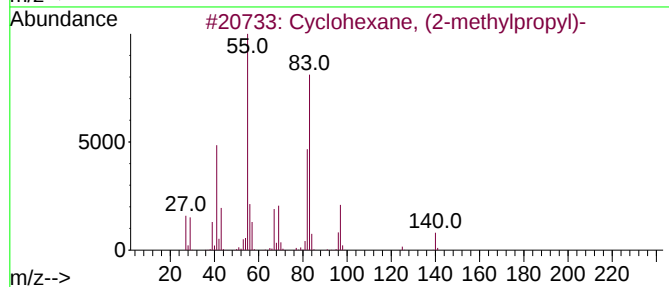
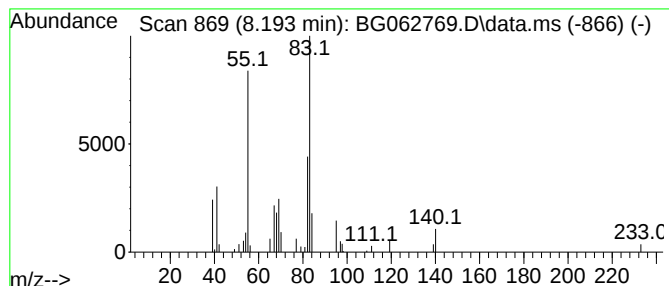
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Cyclic8.193 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.193	3.27 ng/ul	55337	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
2			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	64
4			2-Butyl-.delta.1-pyrroline	125	C8H15N	064319-86-4	50
5			Cyclohexane, butyl-	140	C10H20	001678-93-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

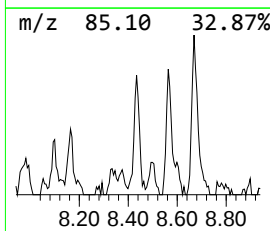
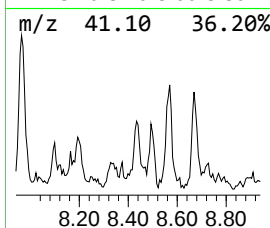
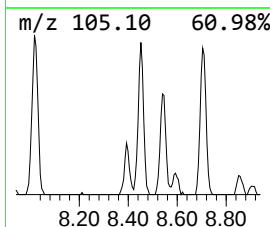
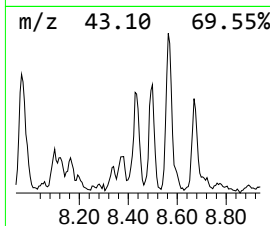
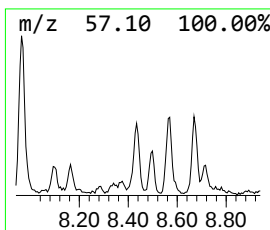
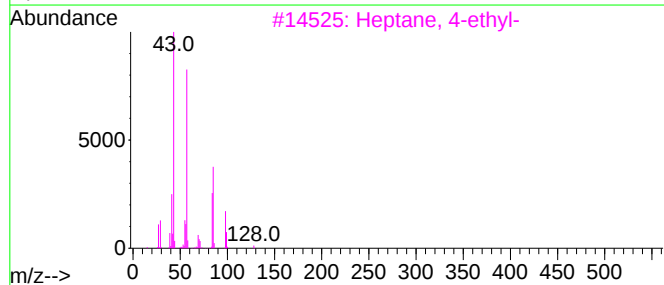
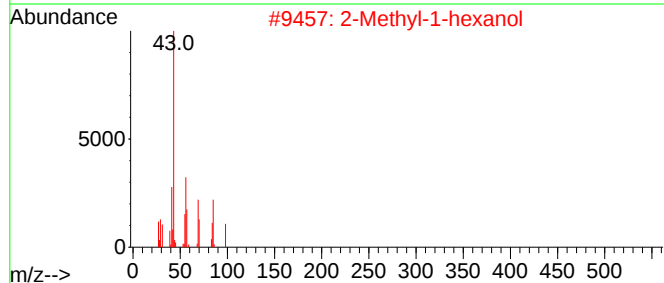
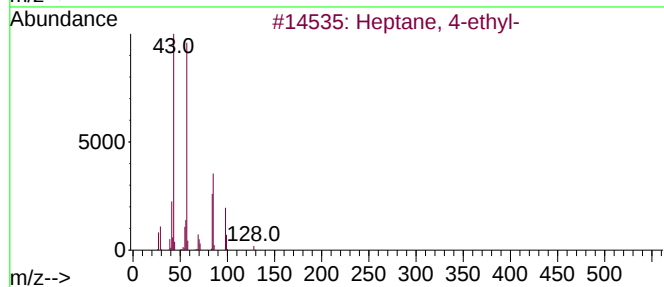
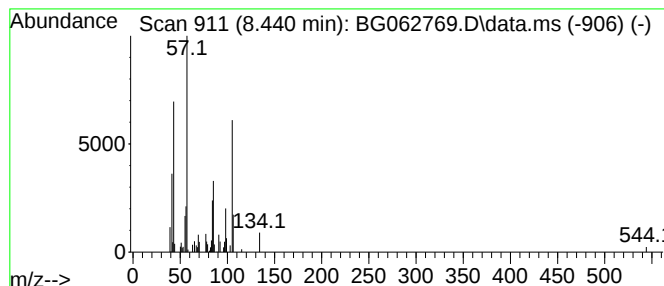
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.440	5.87 ng/ul	99445	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-ethyl-	128	C9H20	002216-32-2	53
2			2-Methyl-1-hexanol	116	C7H16O	1000427-67-0	47
3			Heptane, 4-ethyl-	128	C9H20	002216-32-2	47
4			Heptane, 4-ethyl-	128	C9H20	002216-32-2	43
5			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

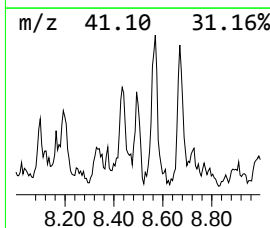
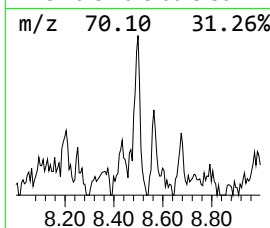
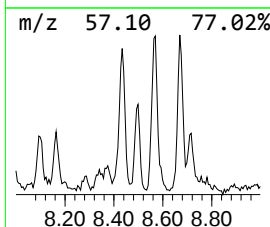
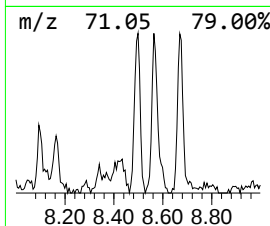
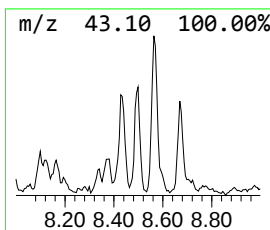
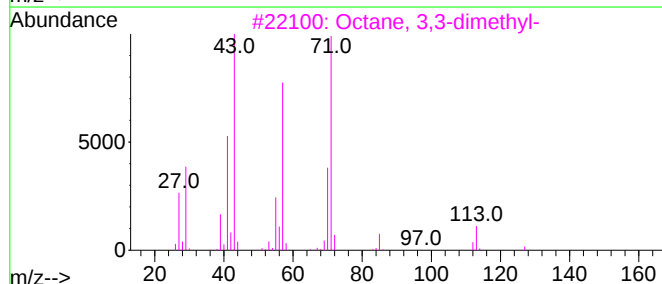
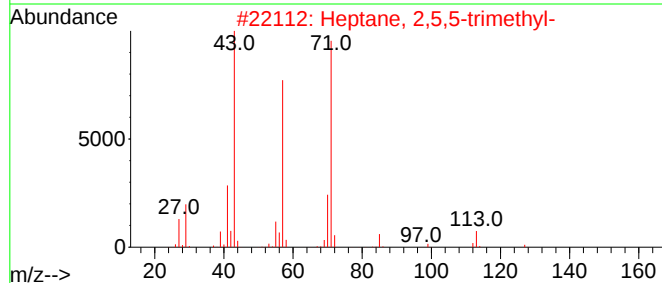
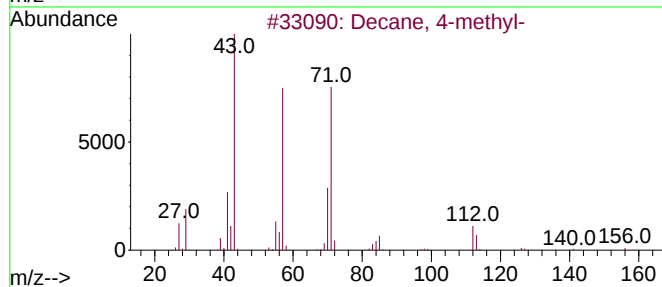
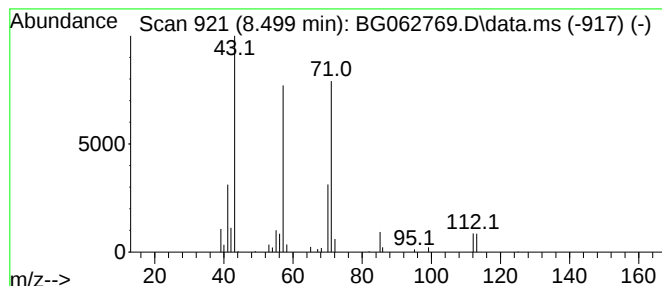
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.499	2.93 ng/ul	49585	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	83
2			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	83
3			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
4			Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	78
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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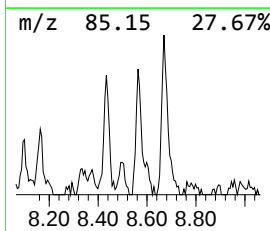
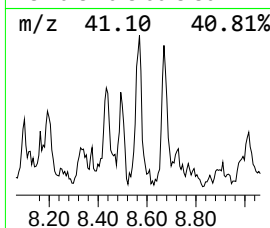
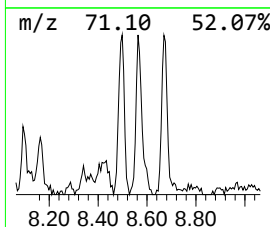
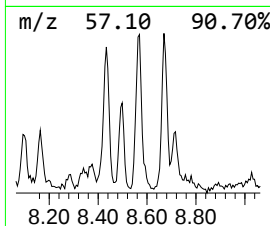
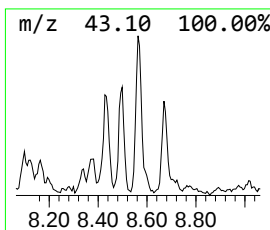
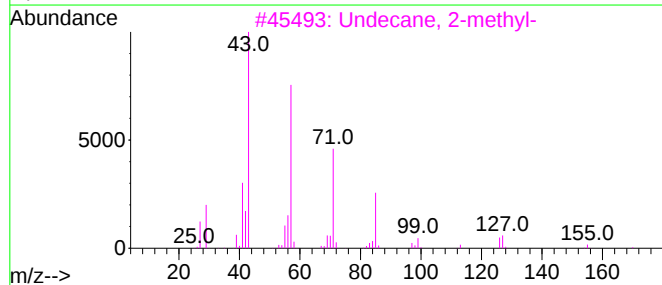
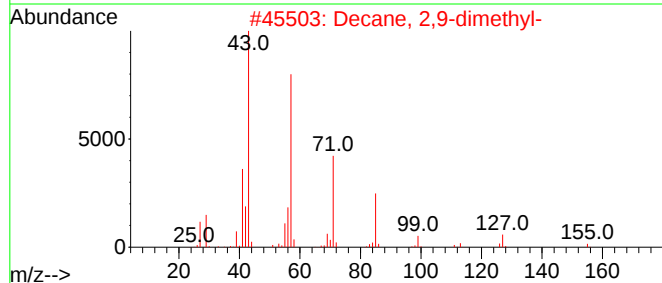
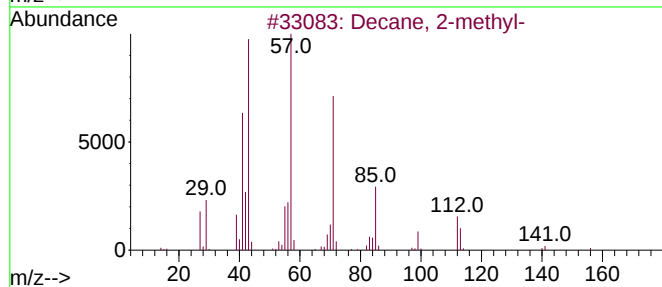
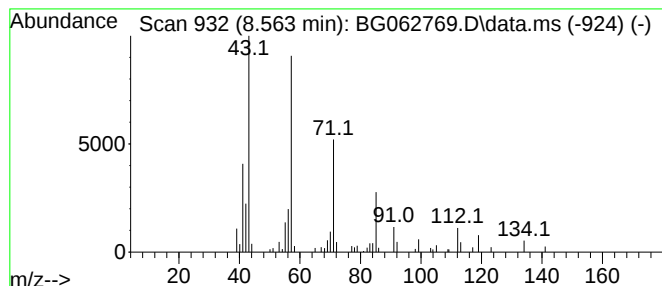
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.563	10.14 ng/ul	171785	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2-methyl-	156	C11H24	006975-98-0	86
2			Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	72
3			Undecane, 2-methyl-	170	C12H26	007045-71-8	64
4			Octane, 2-methyl-	128	C9H20	003221-61-2	64
5			Dodecane, 2-methyl-	184	C13H28	001560-97-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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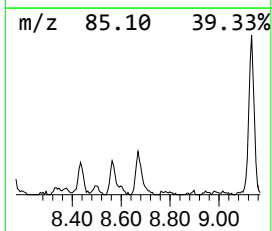
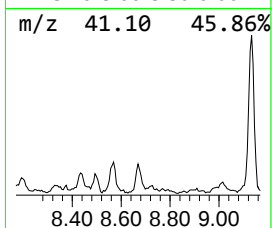
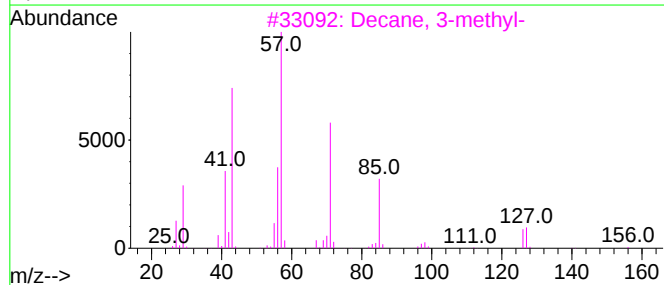
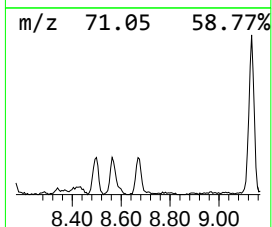
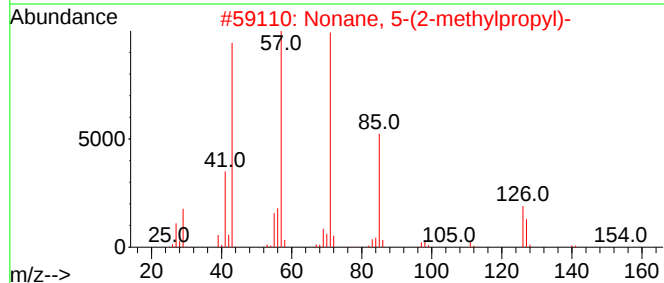
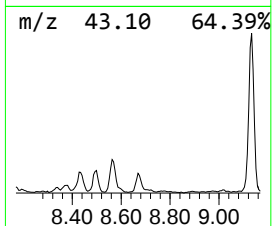
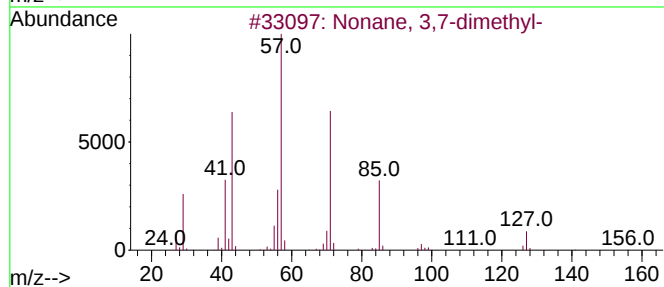
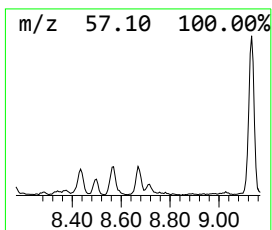
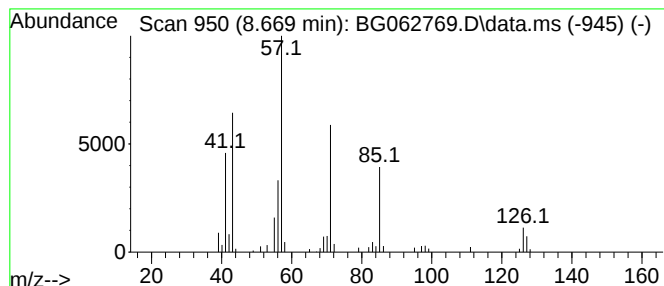
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.669	5.58 ng/ul	94572	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	86
2			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	72
3			Decane, 3-methyl-	156	C11H24	013151-34-3	64
4			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
5			Decane, 3-methyl-	156	C11H24	013151-34-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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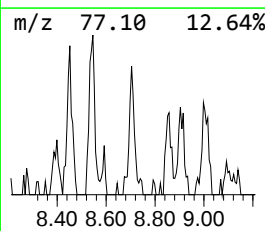
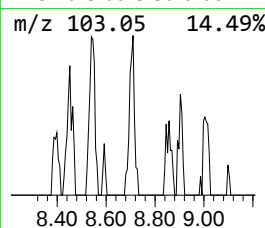
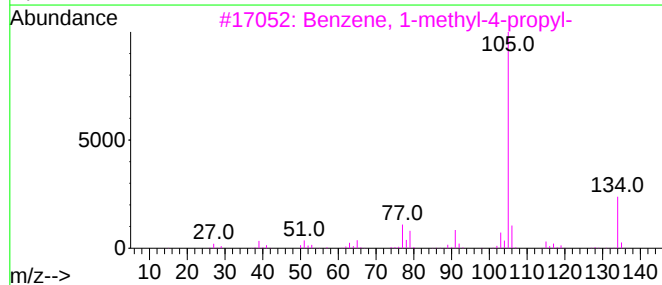
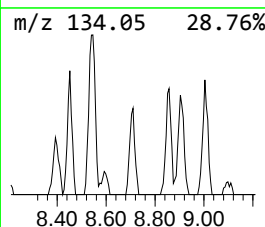
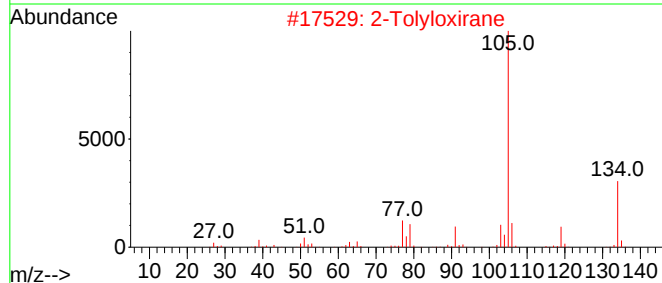
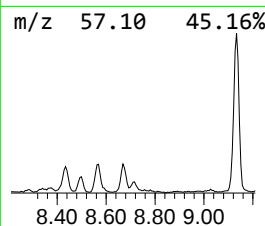
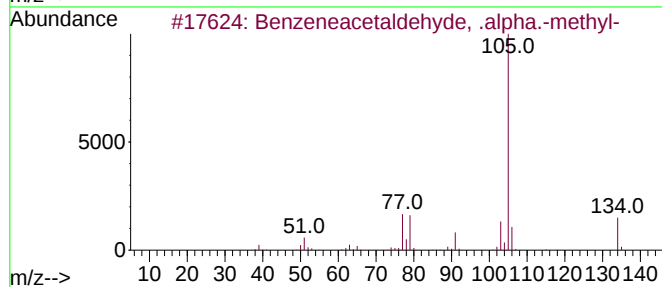
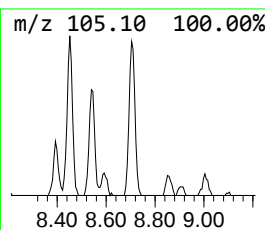
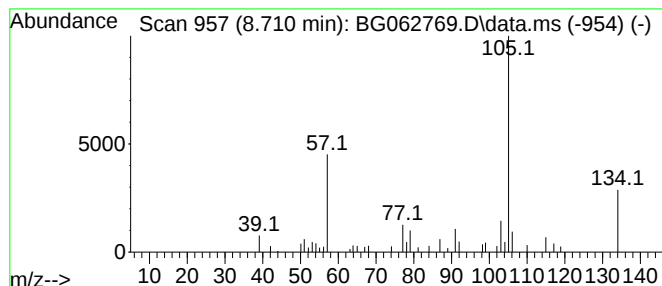
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzeneacetaldehyde, .alpha... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.710	3.74 ng/ul	63370	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	76
2			2-Tolyloxirane	134	C9H10O	002783-26-8	76
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	74
4			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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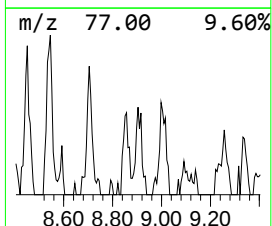
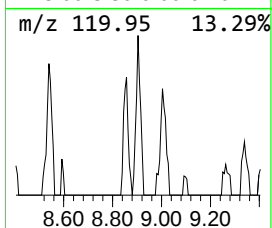
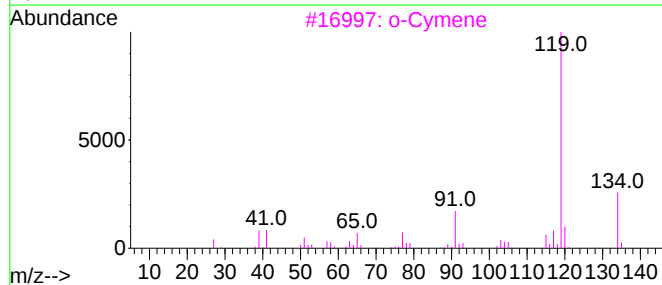
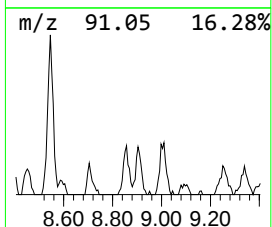
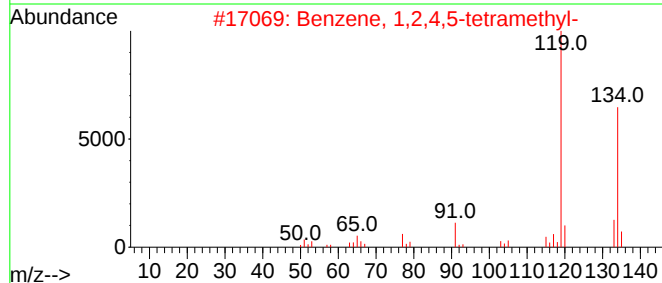
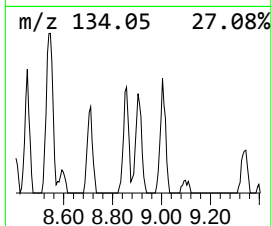
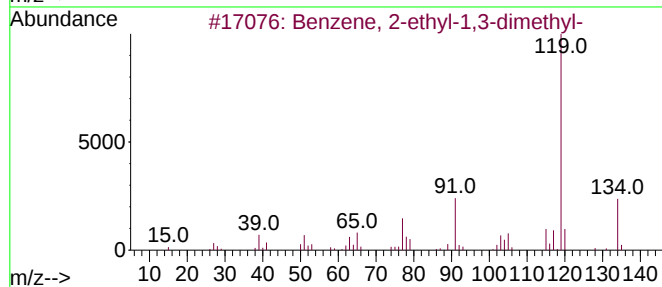
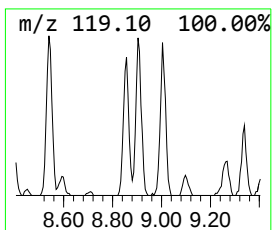
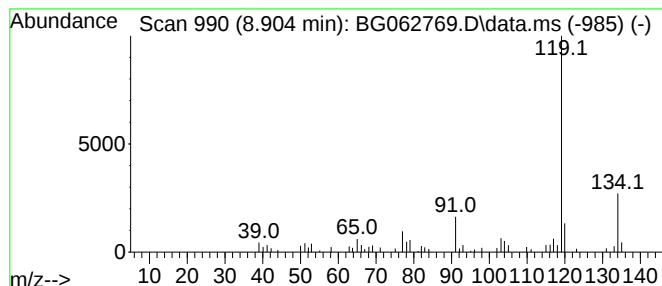
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.904	3.72 ng/ul	63080	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
3			o-Cymene	134	C10H14	000527-84-4	91
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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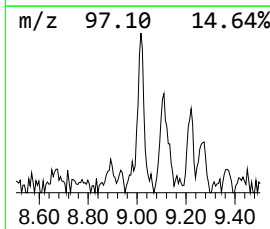
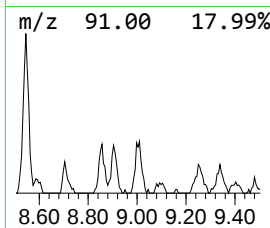
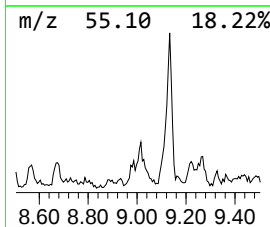
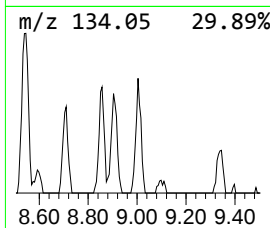
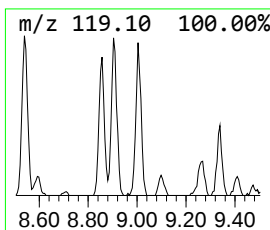
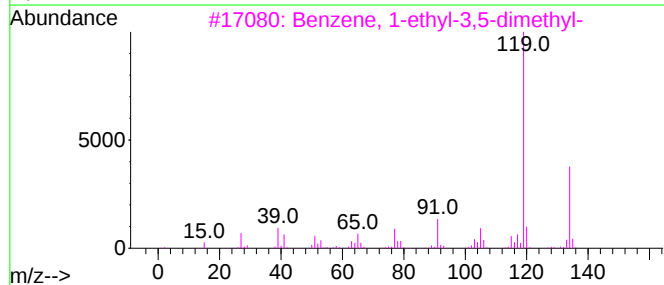
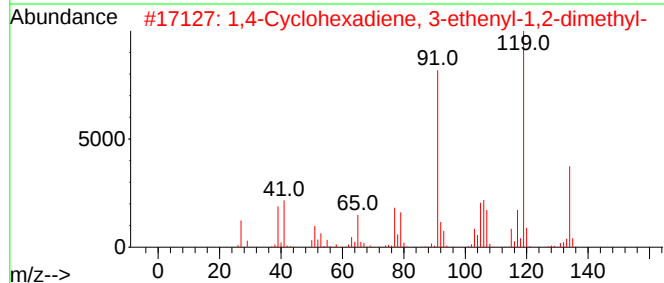
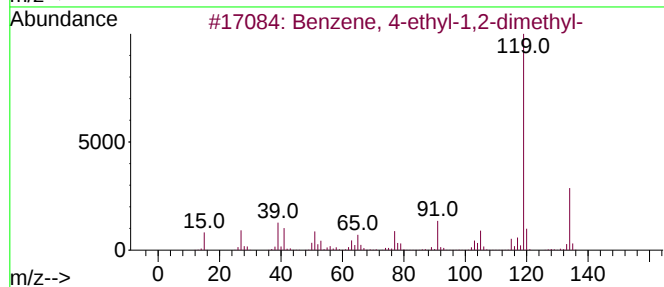
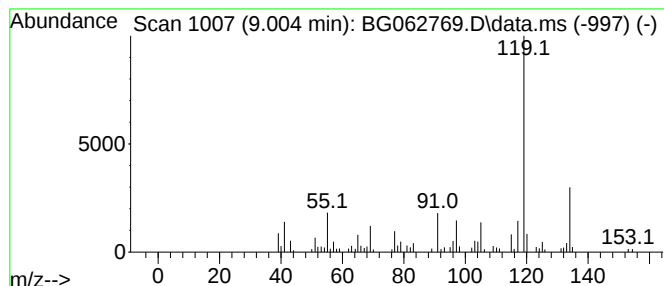
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.004	7.78 ng/ul	131856	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87
2			1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	86
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	83
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

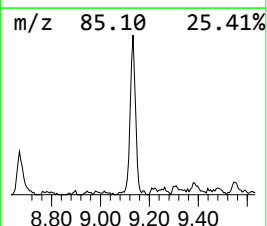
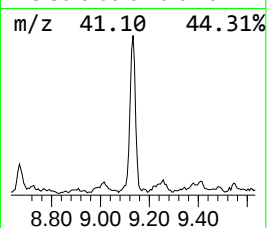
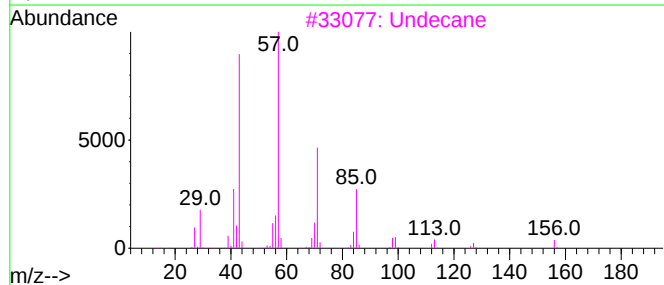
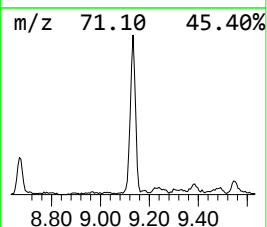
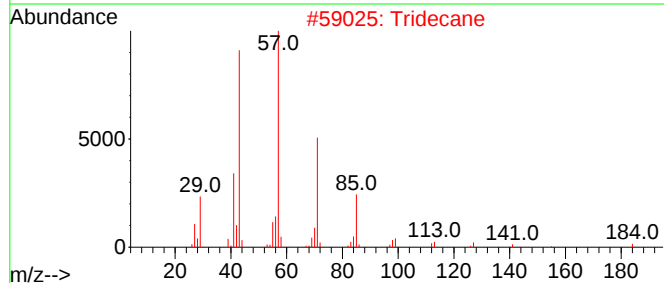
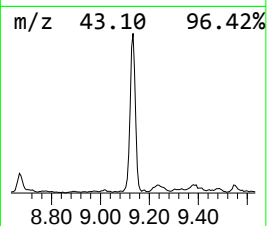
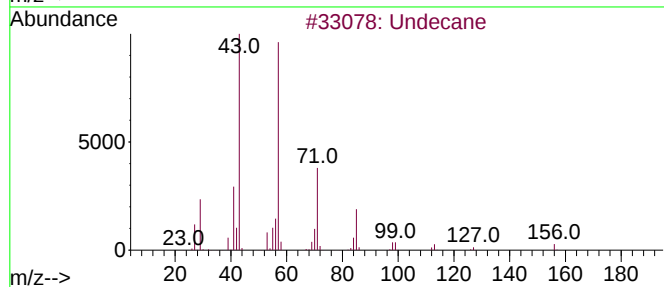
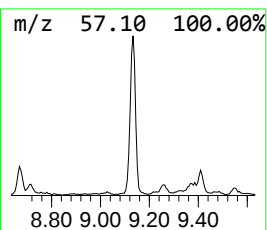
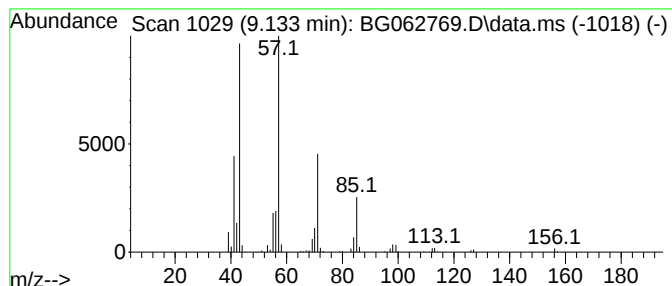
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.133	29.64 ng/ul	502328	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Tridecane	184	C13H28	000629-50-5	86
3			Undecane	156	C11H24	001120-21-4	83
4			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
5			Undecane	156	C11H24	001120-21-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

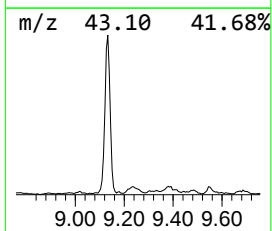
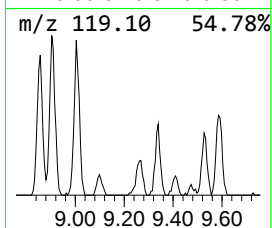
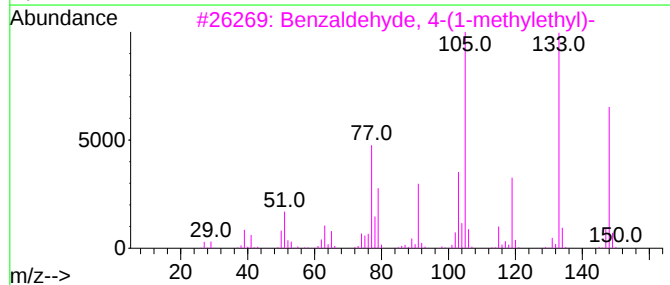
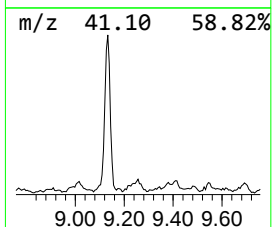
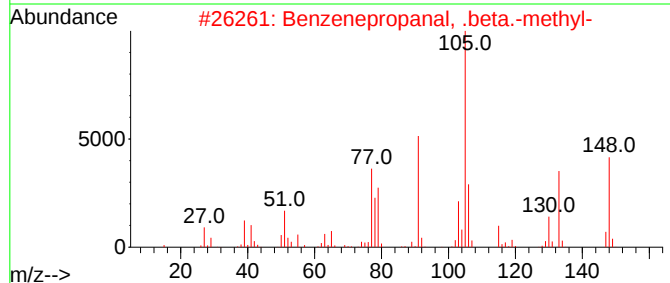
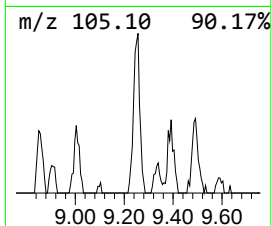
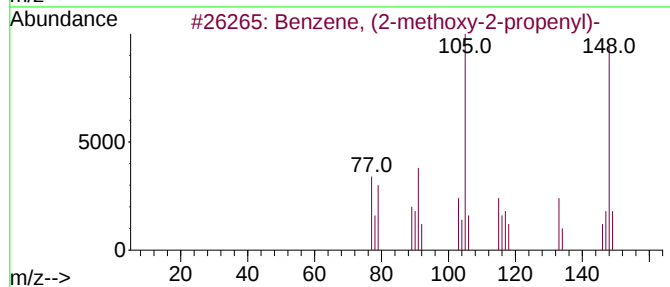
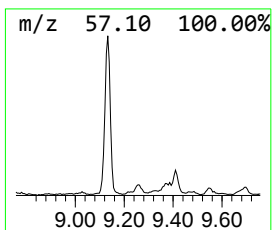
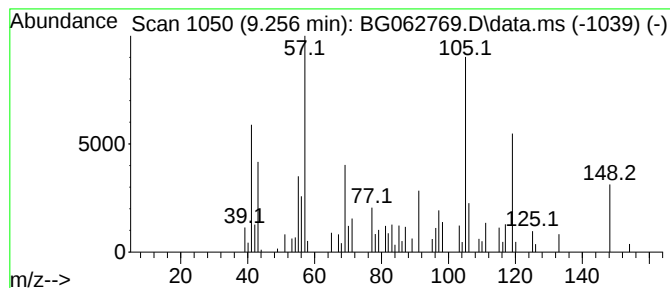
Instrument :
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ClientSampleId :
DCVY4DL2

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 25 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.256	6.59 ng/ul	111623	1,4-Dichlorobenzene-d4		7.905	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, (2-methoxy-2-propenyl)-		148	C10H12O	026473-60-9	25
2	Benzenepropanal, .beta.-methyl-		148	C10H12O	016251-77-7	25
3	Benzaldehyde, 4-(1-methylethyl)-		148	C10H12O	000122-03-2	22
4	Benzene, (1,1-dimethylpropyl)-		148	C11H16	002049-95-8	22
5	2-Butanone, 4-phenyl-		148	C10H12O	002550-26-7	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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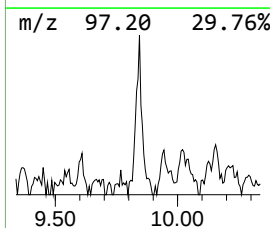
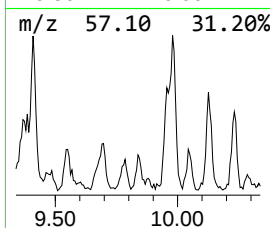
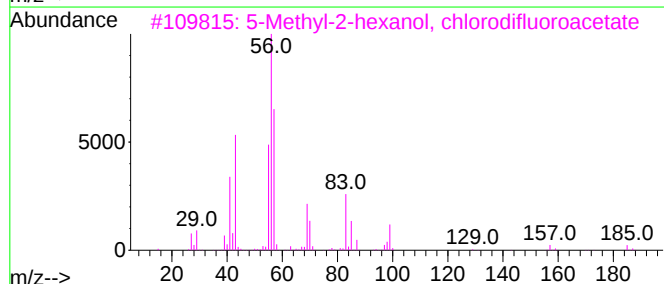
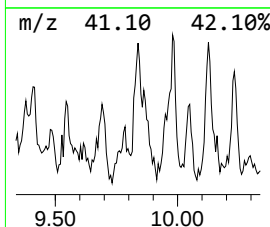
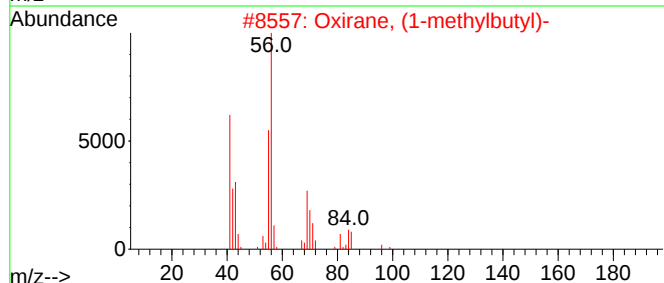
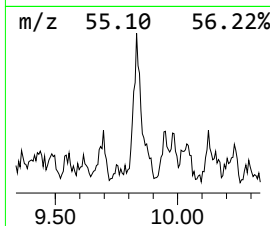
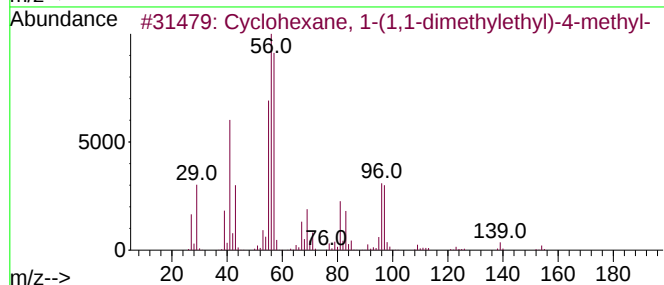
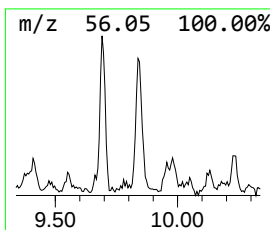
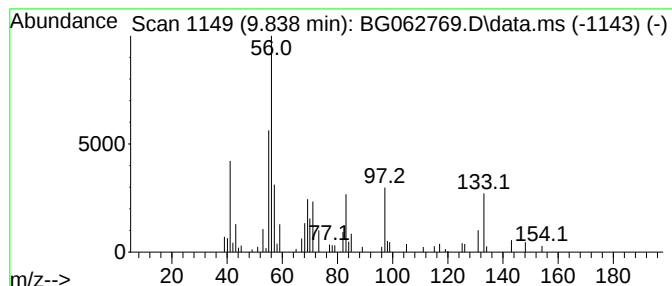
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Cyclic9.838 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.838	2.33 ng/ul	85385	Naphthalene-d8	10.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-(1,1-dimethylethy...	154	C11H22	075736-66-2	49
2			Oxirane, (1-methylbutyl)-	114	C7H14O	053229-39-3	35
3			5-Methyl-2-hexanol, chlorodifluo...	228	C9H15ClF2O2	1010376-27-1	32
4			1-Octene, 2-methyl-	126	C9H18	004588-18-5	30
5			1-Heptene, 2-methyl-	112	C8H16	015870-10-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

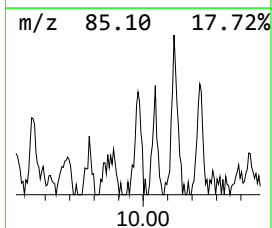
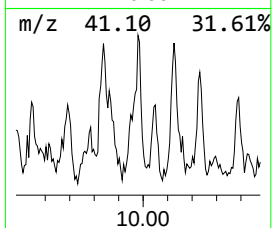
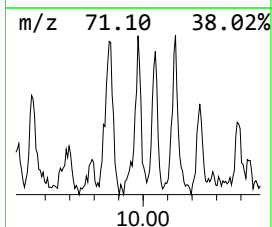
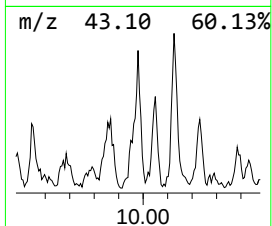
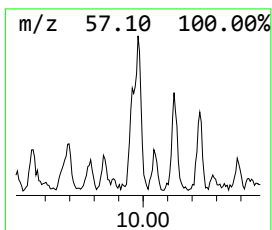
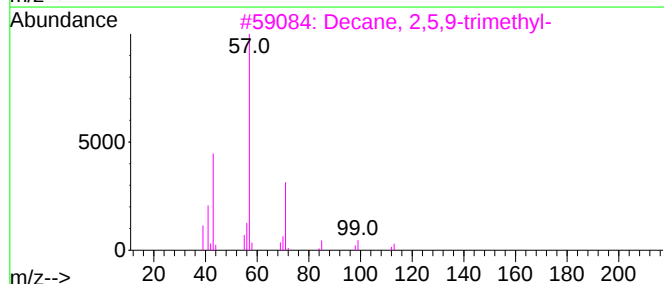
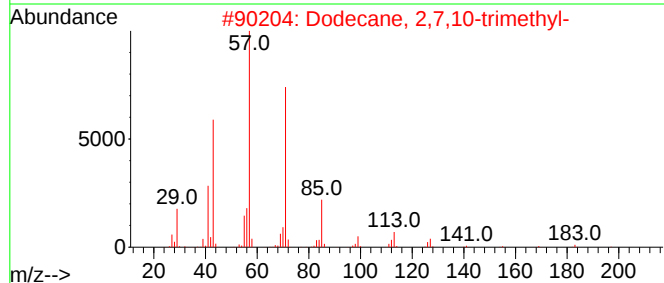
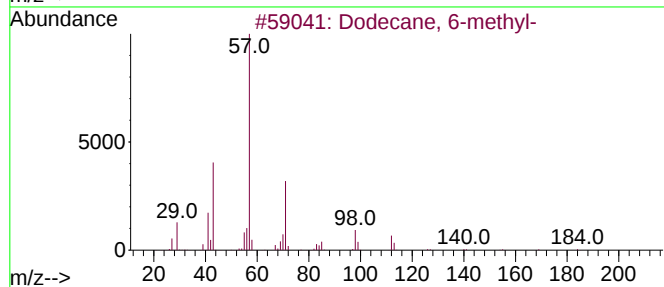
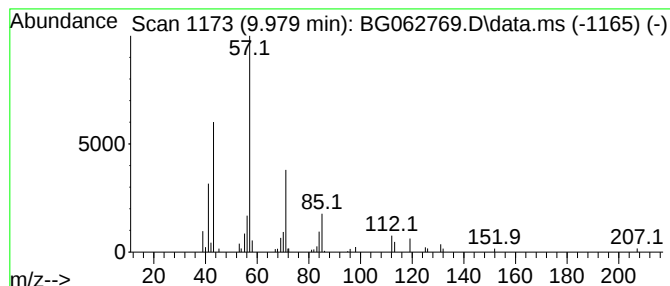
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.979	2.28 ng/ul	83459	Naphthalene-d8	10.696	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 6-methyl-	184	C13H28	006044-71-9	59
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
3	Decane, 2,5,9-trimethyl-	184	C13H28	062108-22-9	53
4	Tridecane, 6-methyl-	198	C14H30	013287-21-3	53
5	Octyl tetradecyl ether	326	C22H46O	1000406-38-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

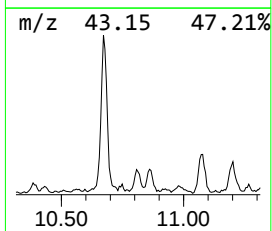
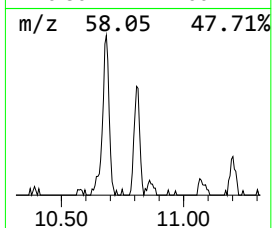
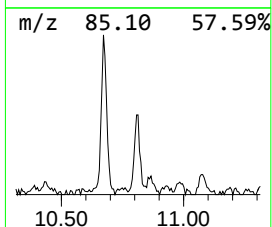
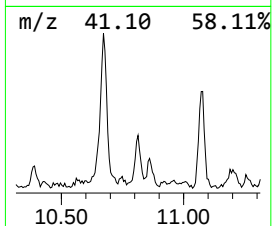
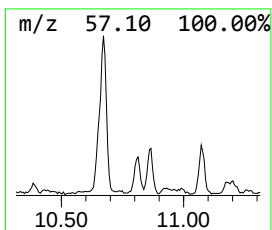
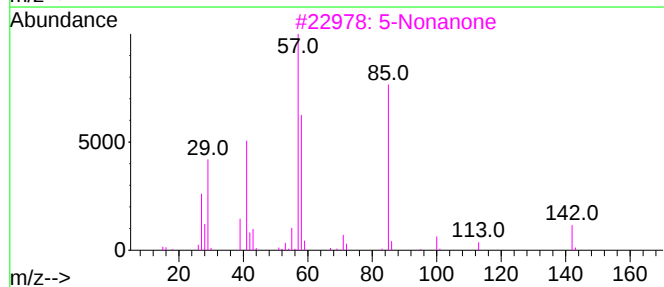
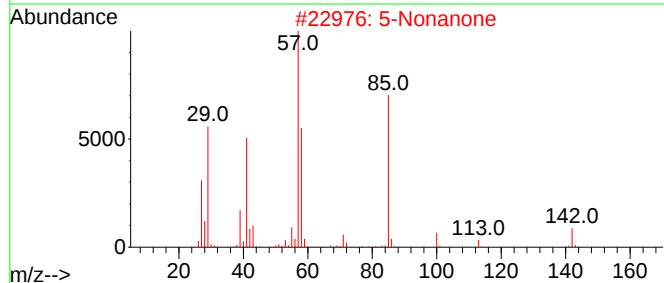
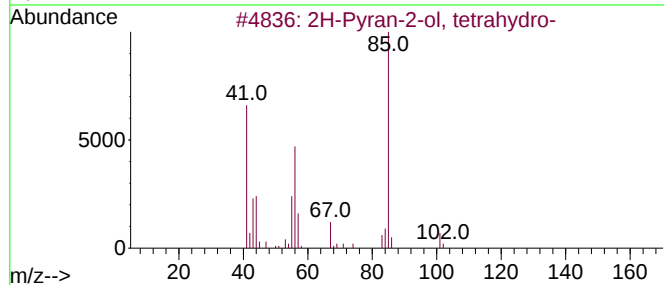
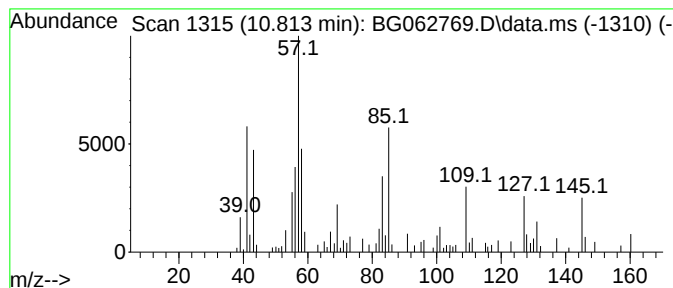
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-02 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.813	2.92 ng/ul	106856	Naphthalene-d8	10.696	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran-2-ol, tetrahydro-	102	C5H10O2	000694-54-2	27
2	5-Nonanone	142	C9H18O	000502-56-7	22
3	5-Nonanone	142	C9H18O	000502-56-7	22
4	Butanoic acid, 3-methyl-, butyl ...	158	C9H18O2	000109-19-3	22
5	1-[(4-Chlorophenyl)sulfonyl]pipe...	260	C10H13ClN2O2S	1000486-20-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

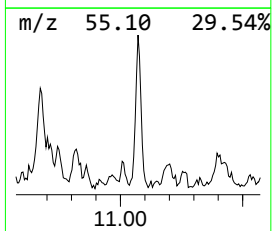
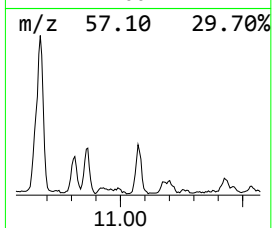
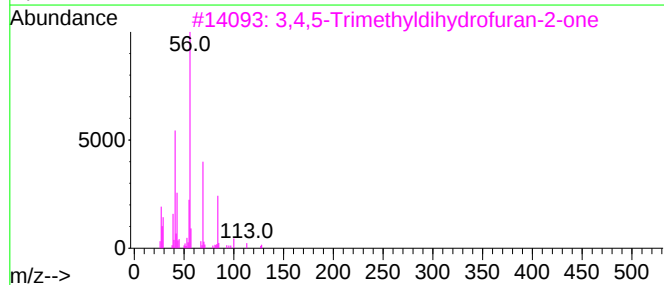
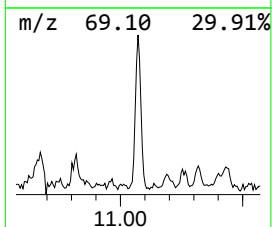
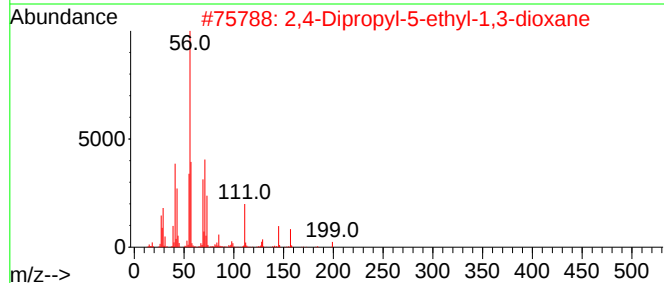
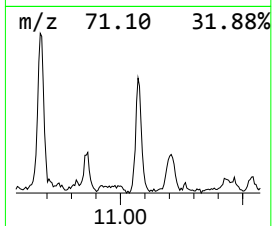
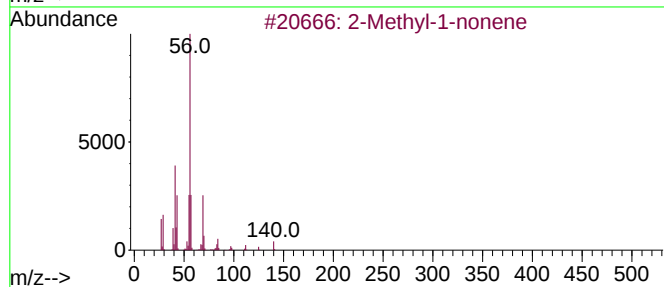
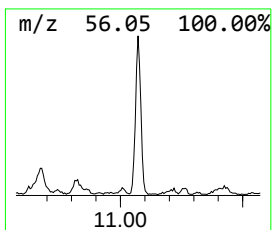
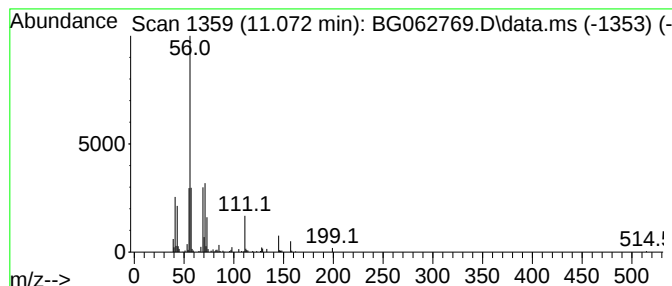
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.072	6.16 ng/ul	225821	Naphthalene-d8	10.696	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-nonene	140	C10H20	002980-71-4	43
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
3	3,4,5-Trimethyldihydrofuran-2-one	128	C7H12O2	1000194-05-2	37
4	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	28
5	Cyclohexylamine	99	C6H13N	000108-91-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

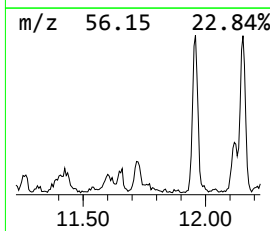
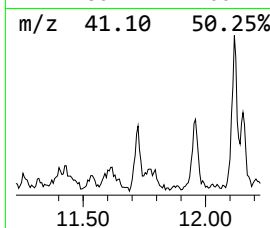
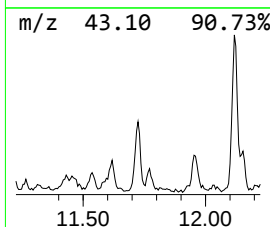
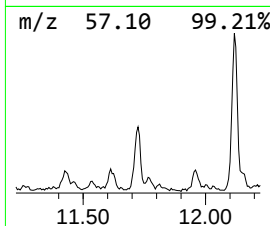
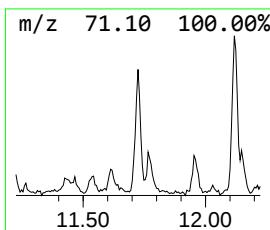
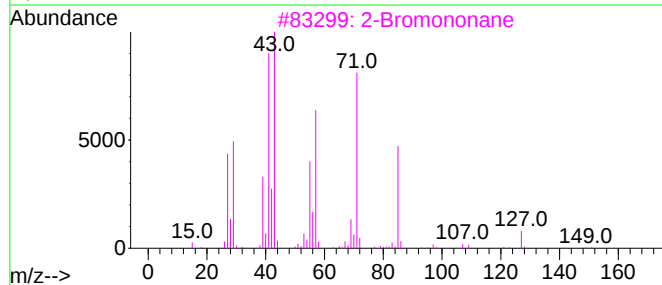
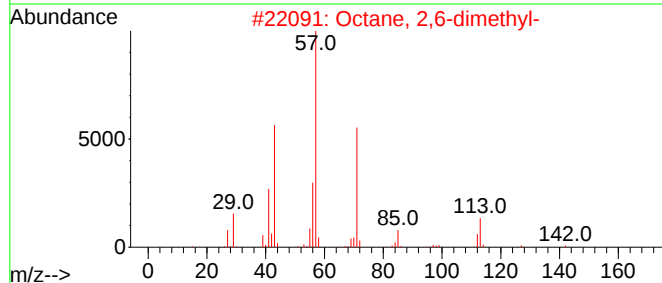
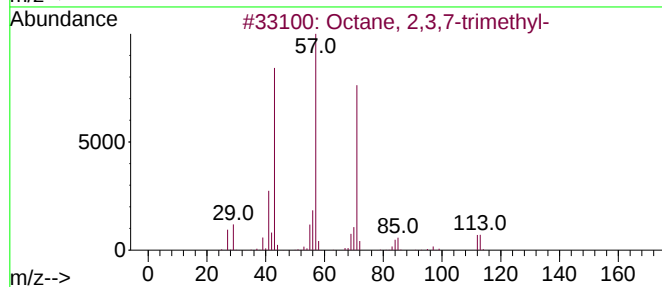
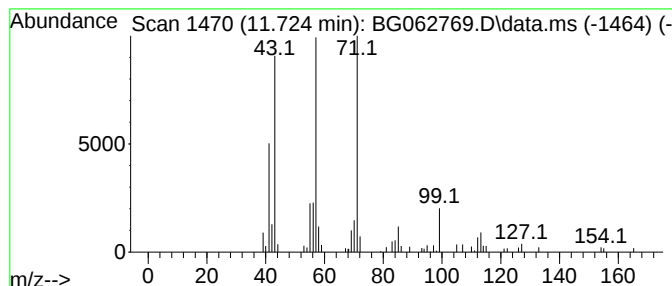
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ClientSampleId :
DCVY4DL2

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Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.724	2.67 ng/ul	97739	Naphthalene-d8	10.696	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3	2-Bromononane	206	C9H19Br	002216-35-5	53
4	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	50
5	4-Heptanone, 3-methyl-	128	C8H16O	015726-15-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

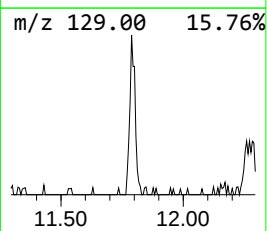
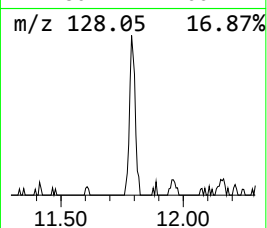
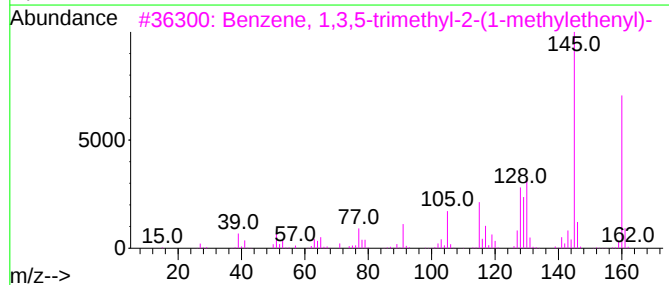
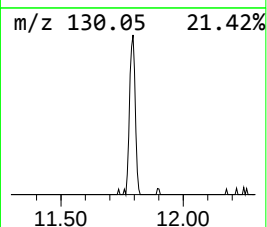
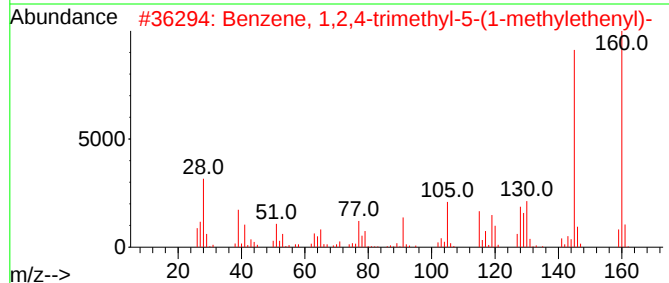
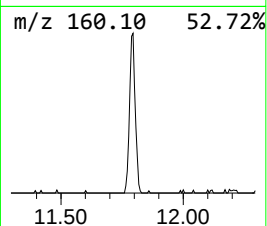
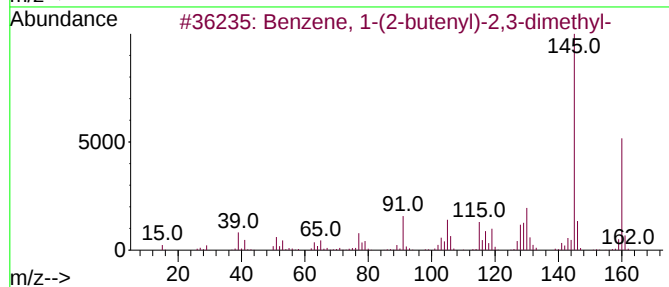
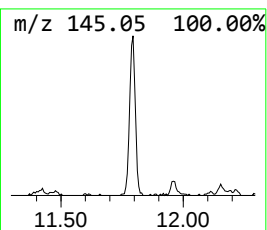
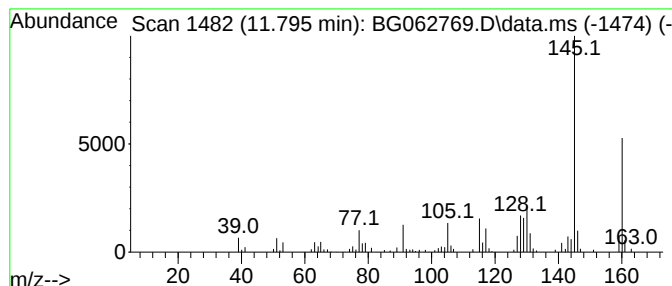
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.795	5.32 ng/ul	194944	Naphthalene-d8	10.696	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	95
2	Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	95
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	95
4	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
5	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

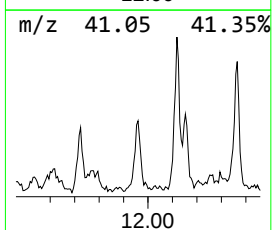
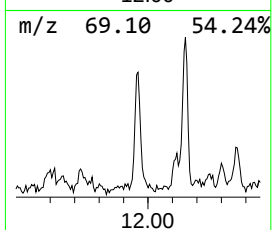
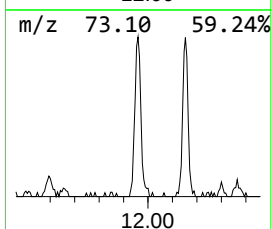
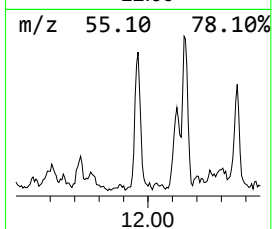
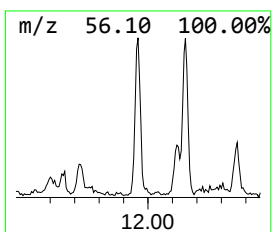
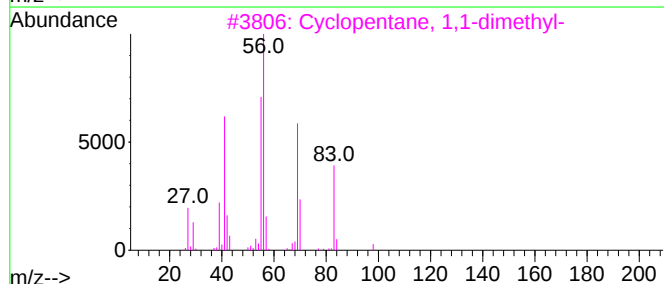
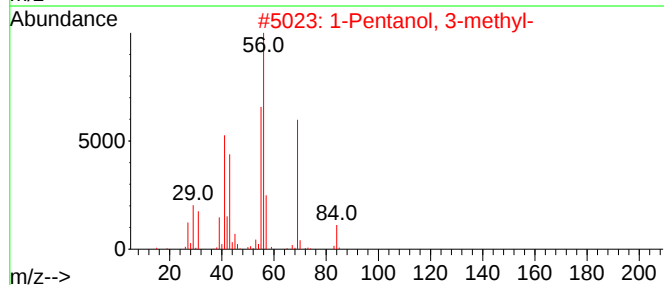
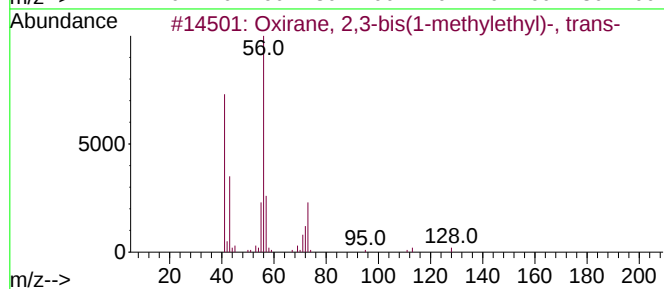
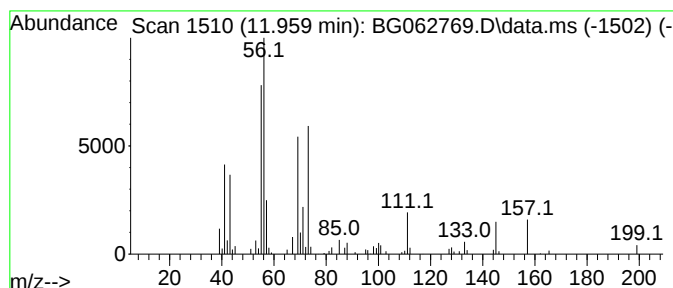
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 unknown-03 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.959	3.51 ng/ul	128672	Naphthalene-d8	10.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	47
2			1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	43
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	43
4			1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	40
5			1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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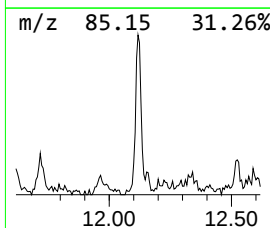
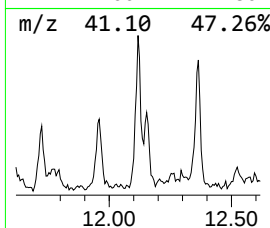
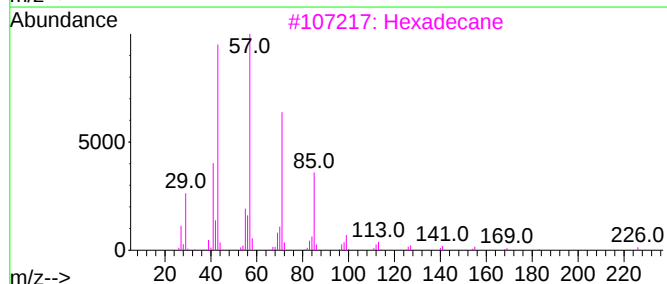
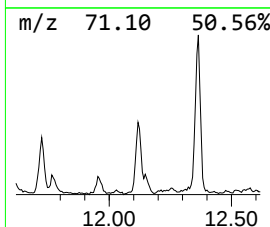
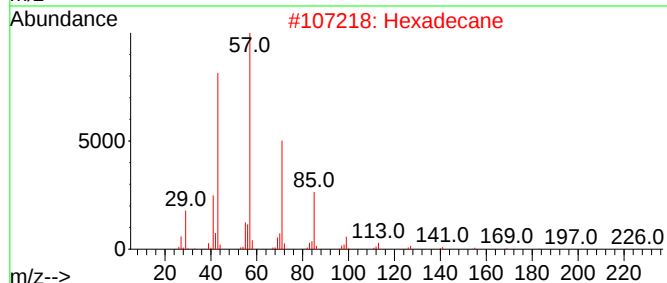
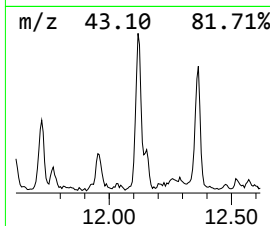
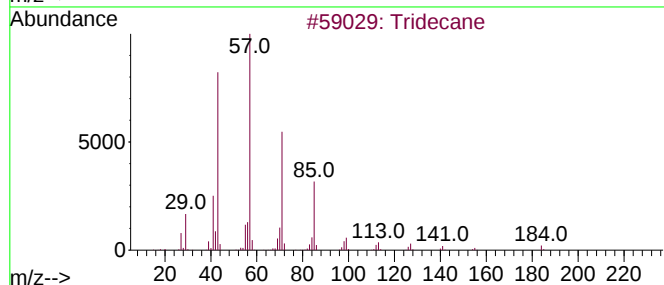
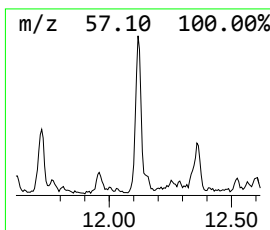
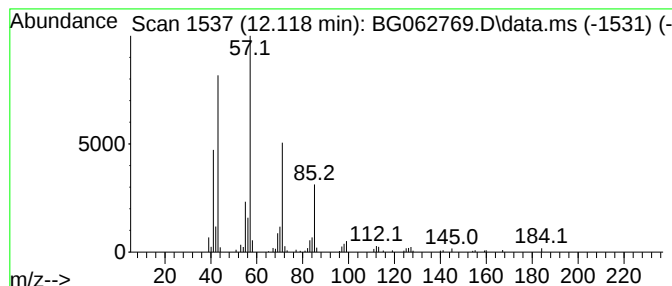
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.118	5.19 ng/ul	190334	Naphthalene-d8	10.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	97
2			Hexadecane	226	C16H34	000544-76-3	90
3			Hexadecane	226	C16H34	000544-76-3	90
4			Tetradecane	198	C14H30	000629-59-4	90
5			Tridecane	184	C13H28	000629-50-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 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

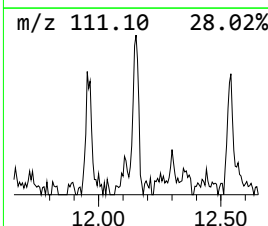
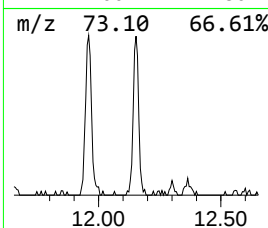
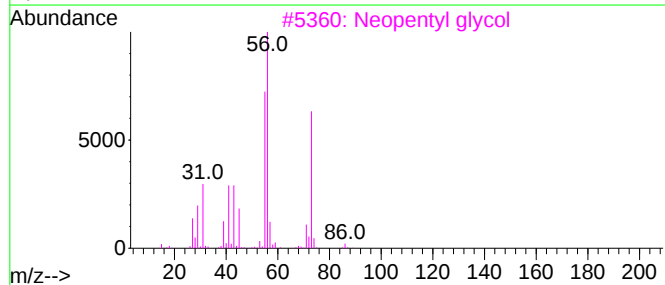
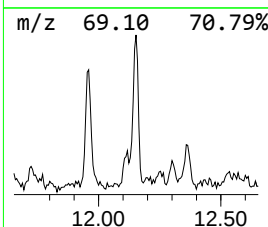
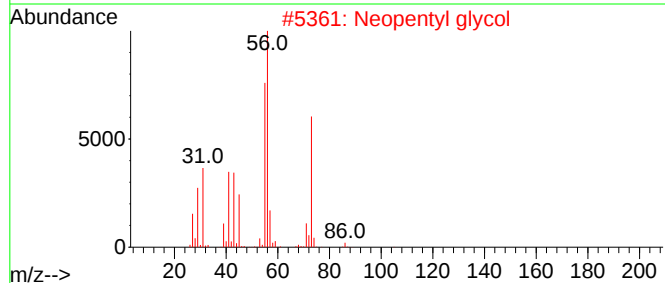
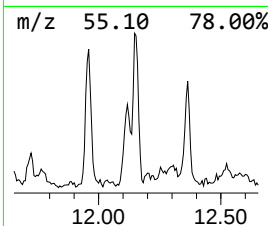
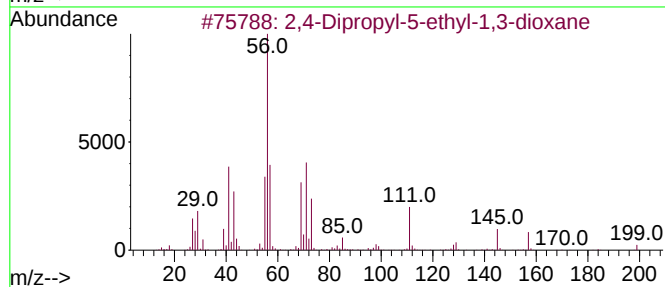
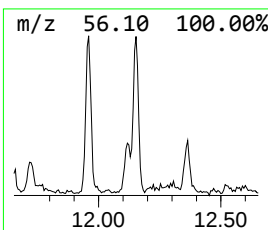
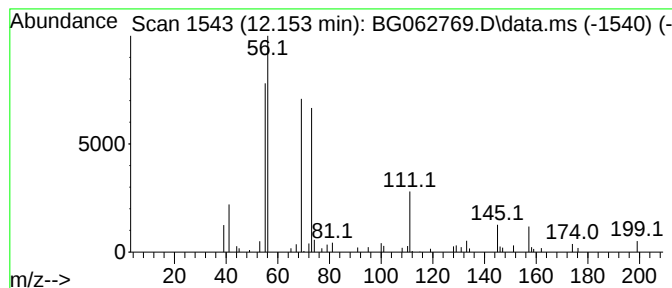
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 unknown-04

Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.153	3.00 ng/ul	109823	Naphthalene-d8	10.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	36
2			Neopentyl glycol	104	C5H12O2	000126-30-7	33
3			Neopentyl glycol	104	C5H12O2	000126-30-7	33
4			3,4,5-Trimethyldihydrofuran-2-one	128	C7H12O2	1000194-05-2	32
5			Hexyl trifluoroacetate	198	C8H13F3O2	000400-61-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

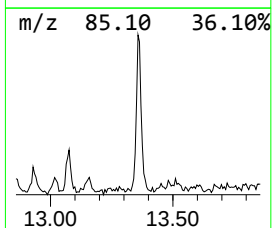
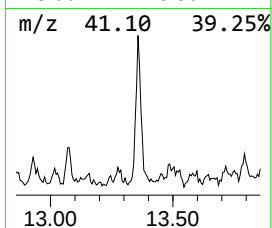
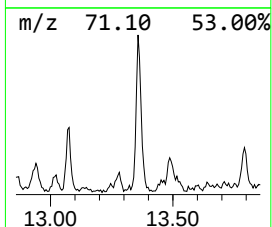
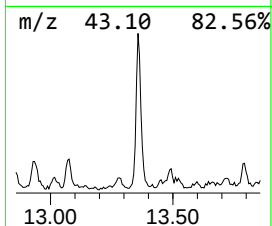
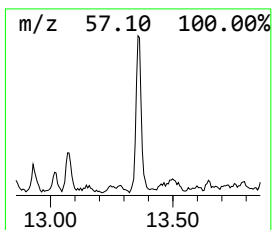
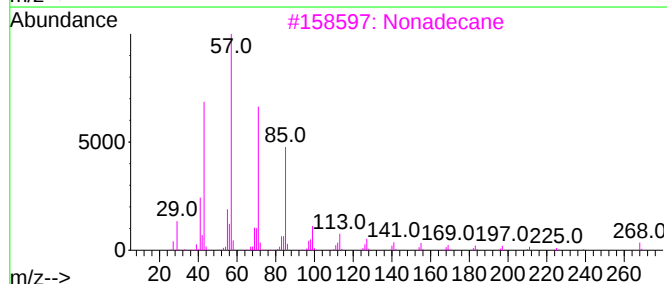
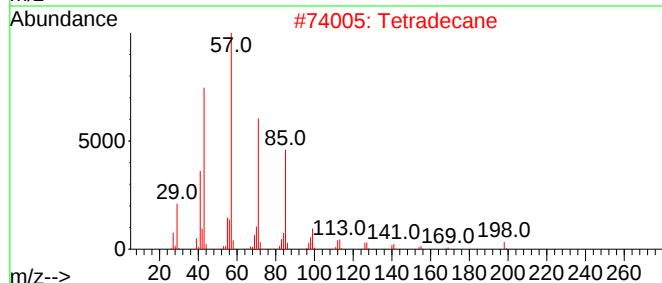
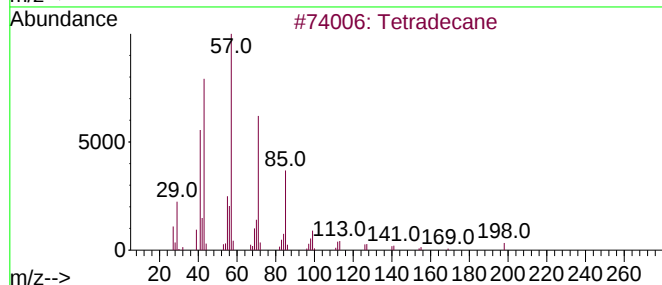
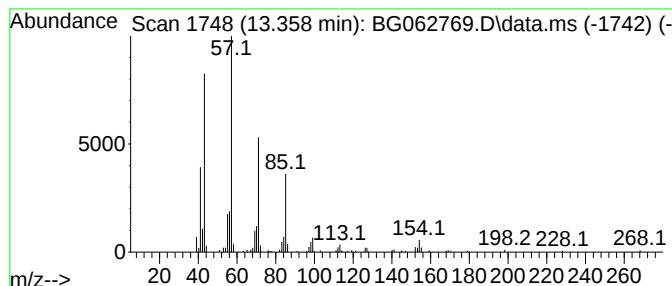
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.358	5.04 ng/ul	223346	Acenaphthene-d10		14.533	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	93
2	Tetradecane		198	C14H30	000629-59-4	93
3	Nonadecane		268	C19H40	000629-92-5	91
4	Nonadecane		268	C19H40	000629-92-5	91
5	Decane, 1-iodo-		268	C10H21I	002050-77-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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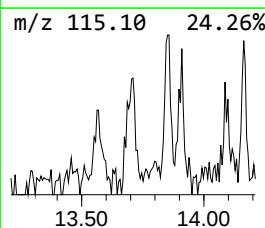
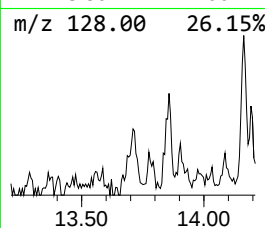
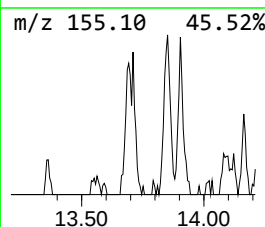
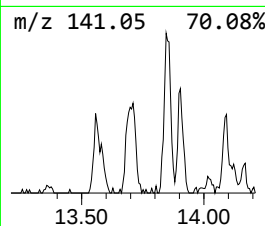
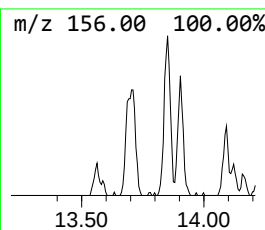
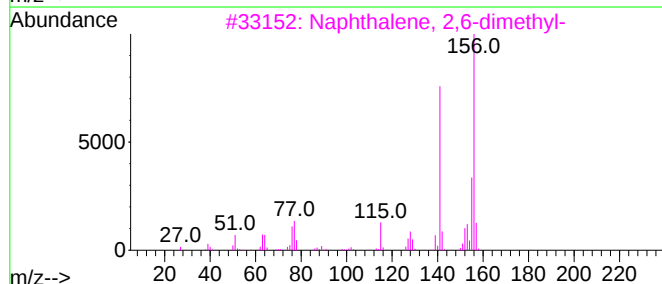
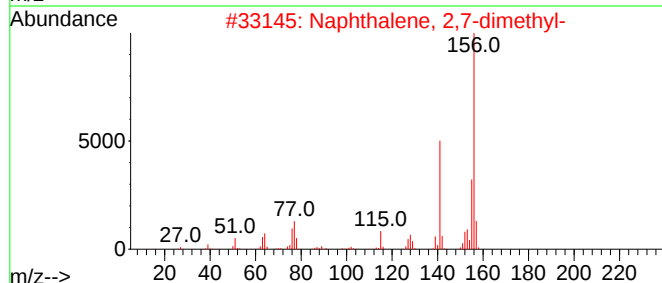
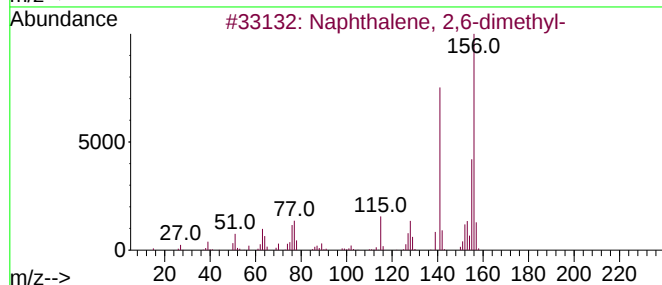
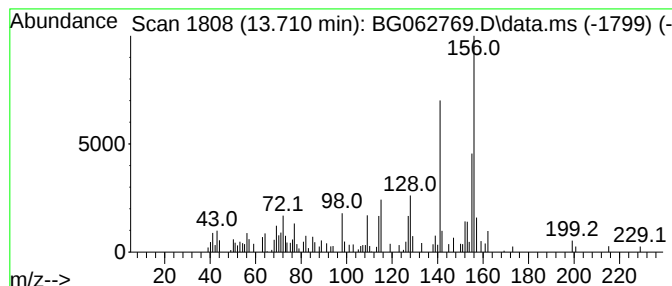
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Naphthalene, 2,6-dimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.710	2.49 ng/ul	110389	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	90
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	90
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	81
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

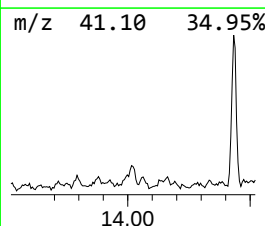
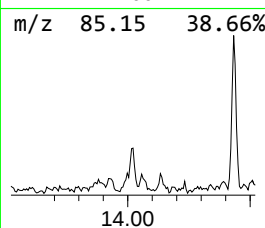
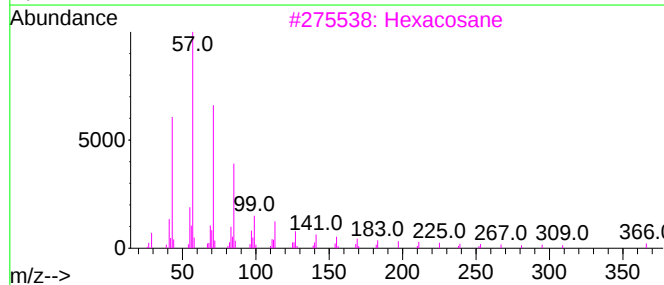
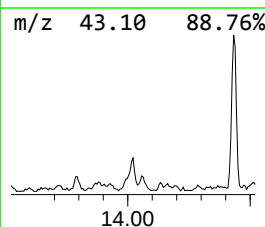
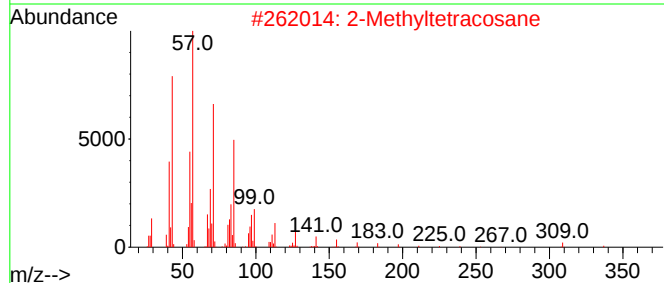
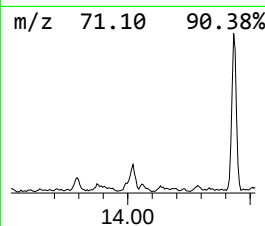
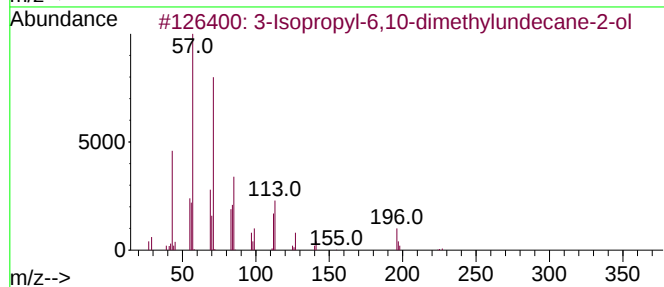
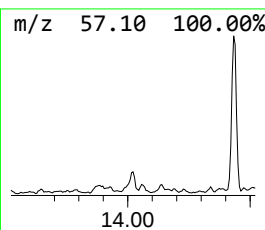
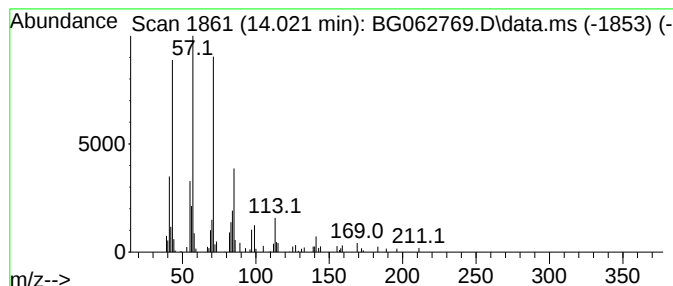
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 3-Isopropyl-6,10-dimethylun... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.022	2.25 ng/ul	99709	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Isopropyl-6,10-dimethylundecan...	242	C16H34O	1000432-47-1	78
2			2-Methyltetracosane	352	C25H52	001560-78-7	58
3			Hexacosane	366	C26H54	000630-01-3	50
4			Hexadecane	226	C16H34	000544-76-3	50
5			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

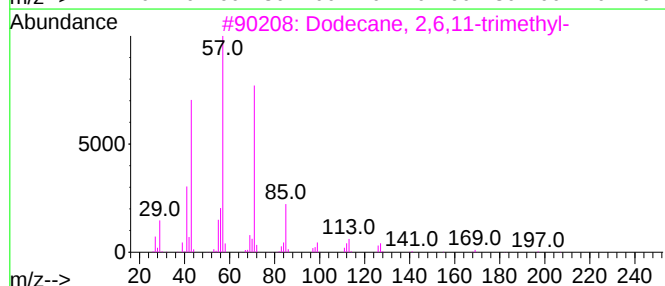
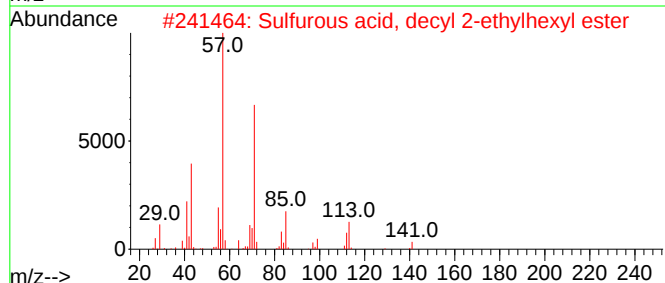
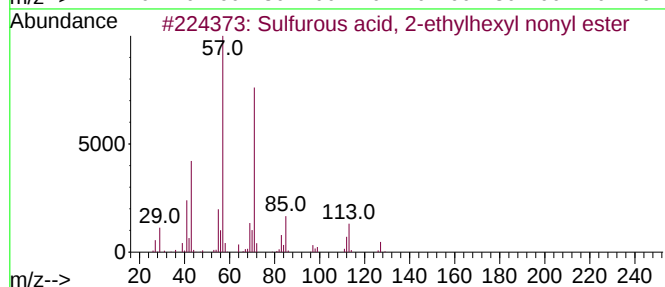
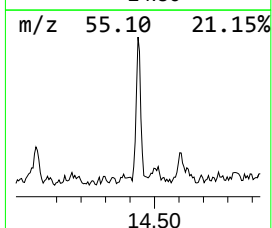
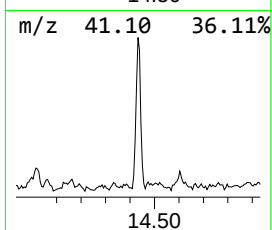
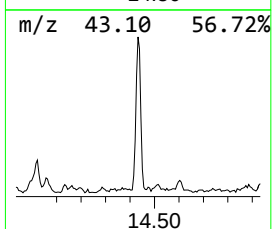
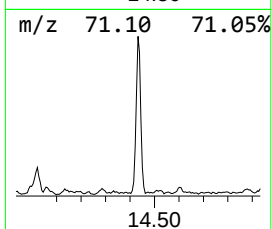
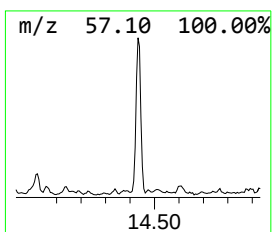
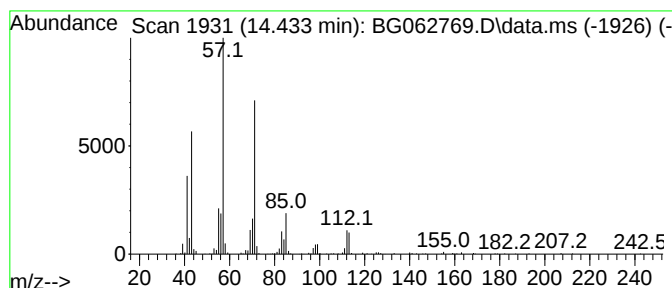
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 Sulfurous acid, 2-ethylhexy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.433	10.15 ng/ul	450141	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	86
2			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	86
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
4			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	72
5			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

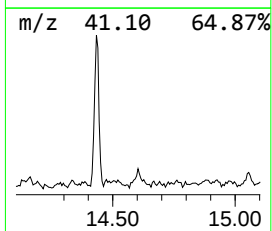
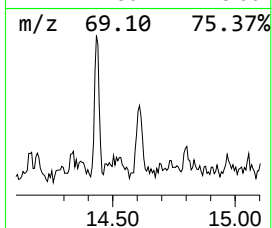
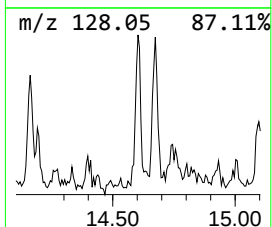
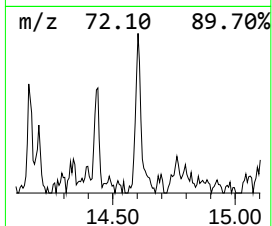
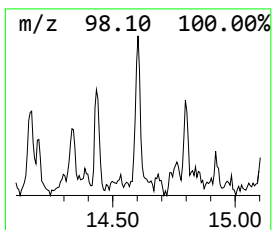
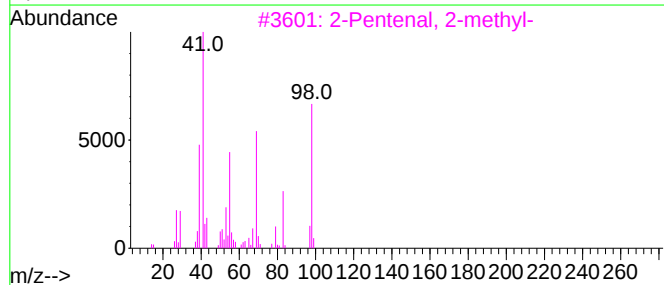
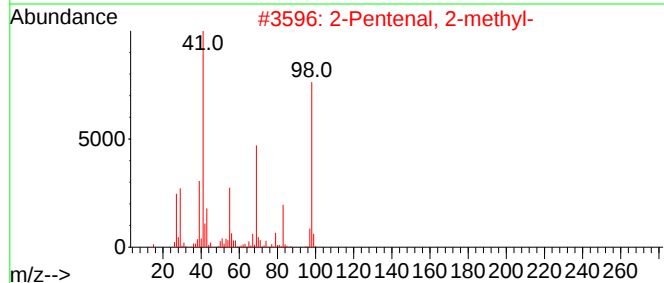
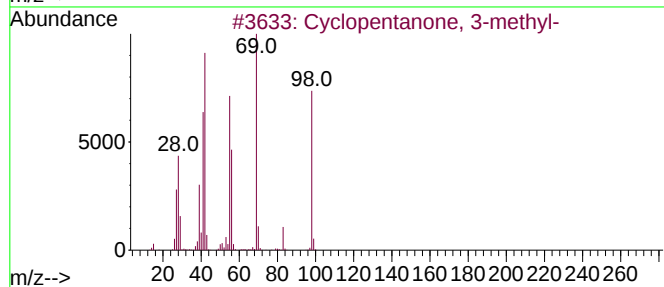
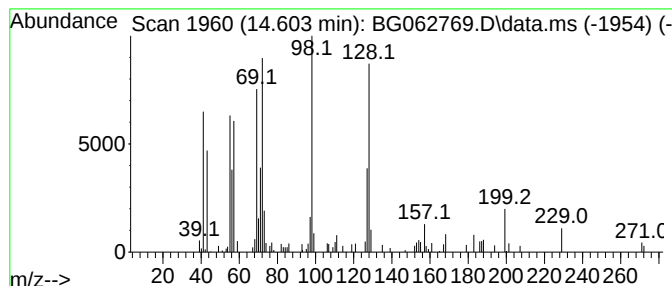
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 unknown-05 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.603	2.08 ng/ul	92287	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone, 3-methyl-	98	C6H10O	001757-42-2	27
2			2-Pentenal, 2-methyl-	98	C6H10O	000623-36-9	14
3			2-Pentenal, 2-methyl-	98	C6H10O	000623-36-9	14
4			2-Ethyl-4-methylthiazole	127	C6H9NS	015679-12-6	14
5			2-Pentenal, 2-methyl-	98	C6H10O	000623-36-9	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

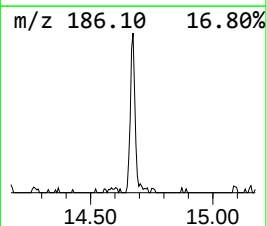
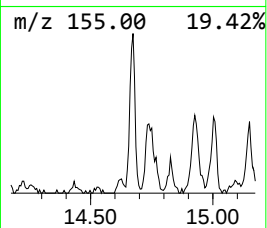
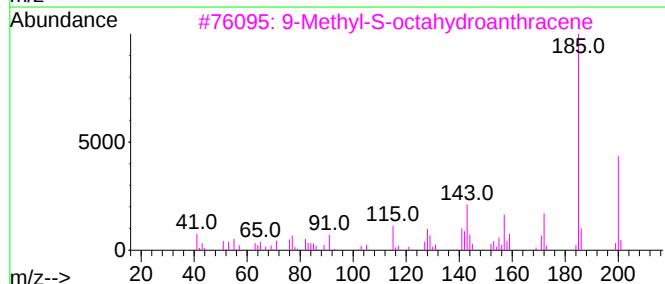
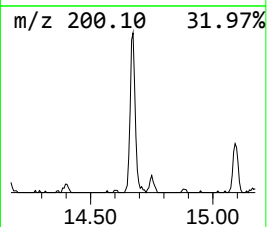
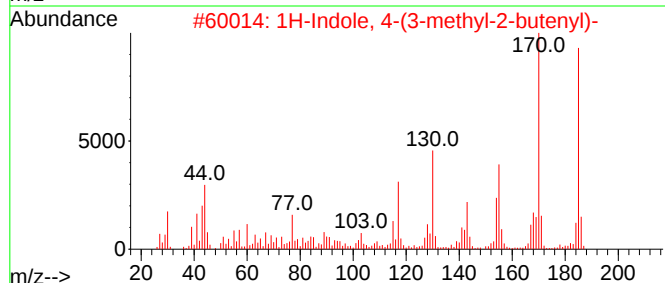
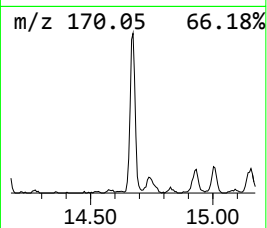
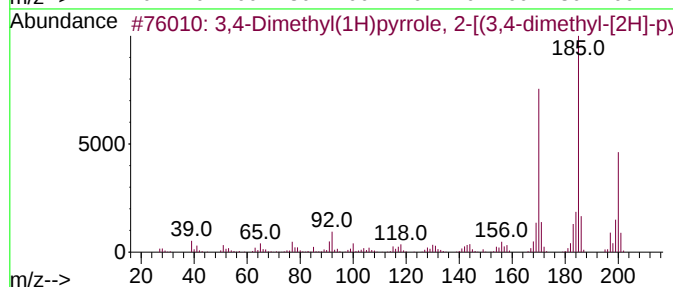
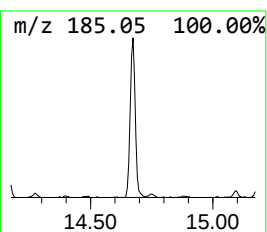
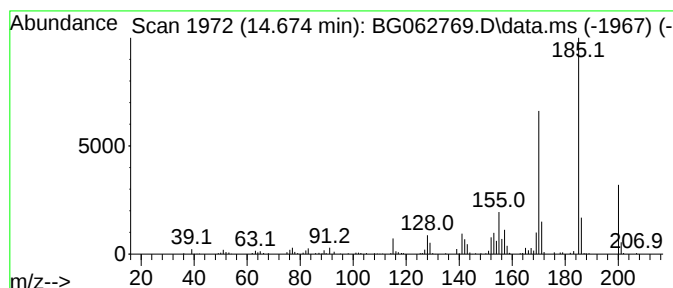
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 40 3,4-Dimethyl(1H)pyrrole, 2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.674	6.79 ng/ul	301158	Acenaphthene-d10	14.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethyl(1H)pyrrole, 2-[(3,4...	200	C13H16N2	065539-79-9	72
2	1H-Indole, 4-(3-methyl-2-butenyl)-	185	C13H15N	032962-25-7	72
3	9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	53
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	52
5	4'-(1-Pyrroyl)acetophenone	185	C12H11NO	022106-37-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

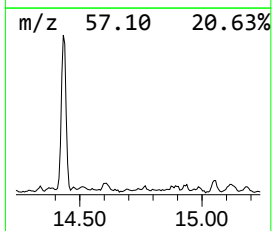
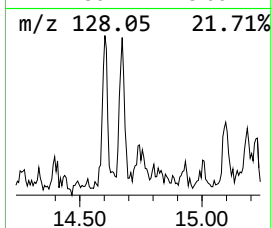
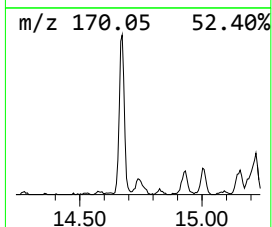
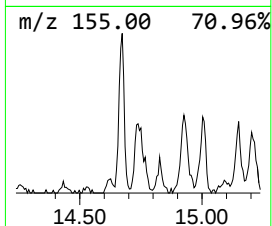
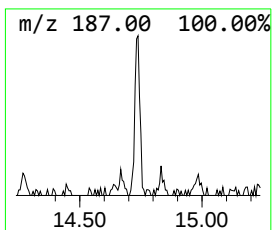
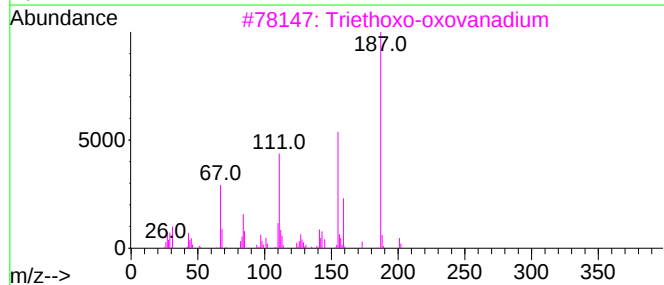
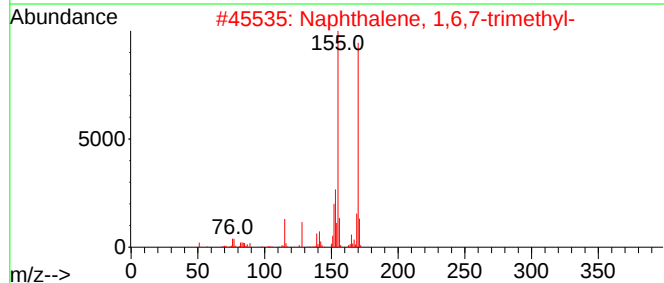
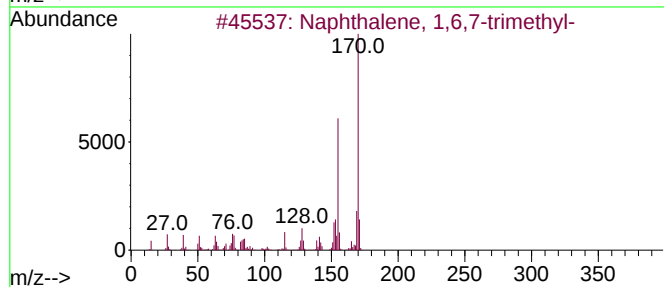
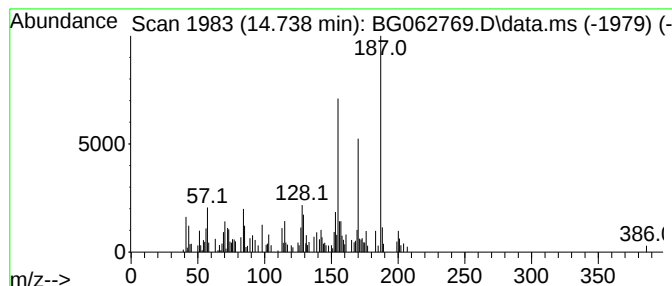
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Naphthalene, 1,6,7-trimethyl- Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.738	2.14 ng/ul	95064	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	59
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	55
3			Triethoxo-oxovanadium	202	C6H15O4V	001686-22-2	49
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	42
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

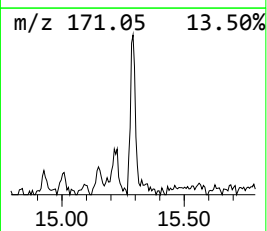
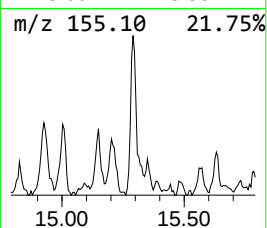
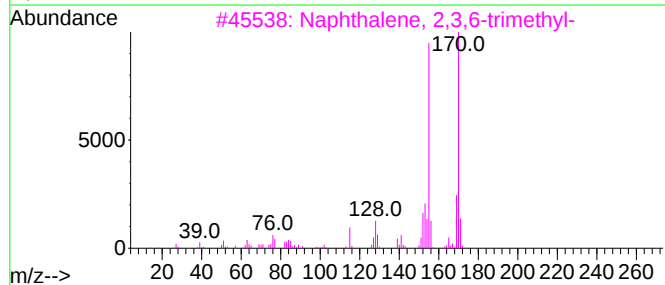
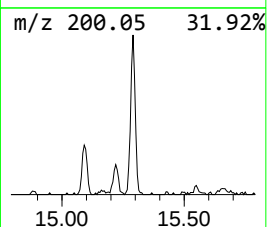
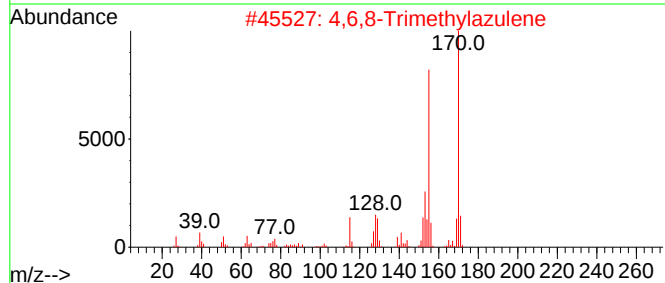
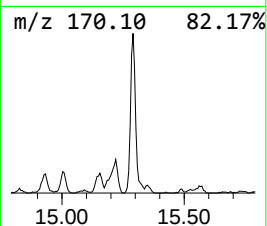
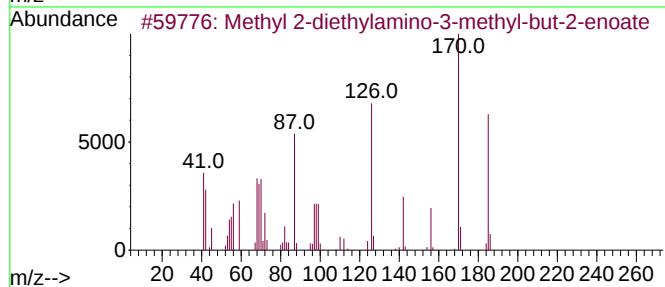
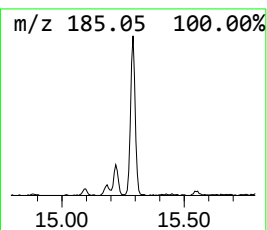
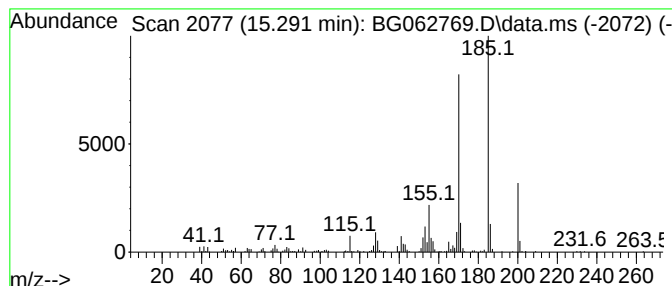
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Methyl 2-diethylamino-3-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.291	7.36 ng/ul	326189	Acenaphthene-d10	14.533

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	52
2		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	43
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	43
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	41
5		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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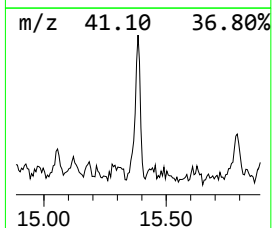
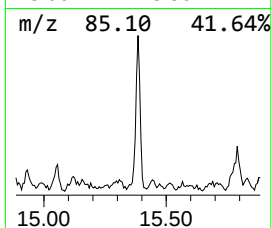
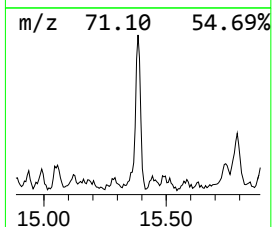
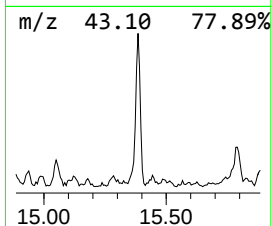
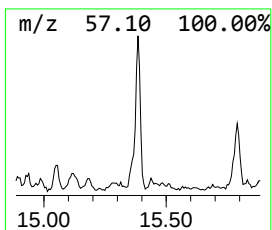
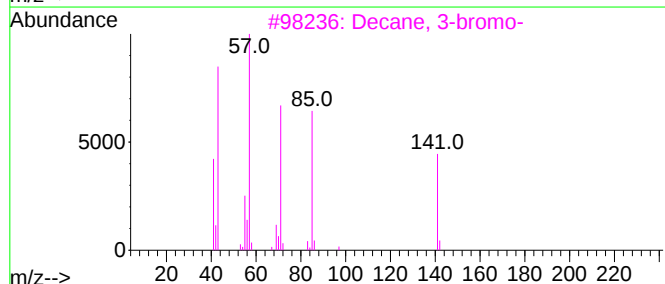
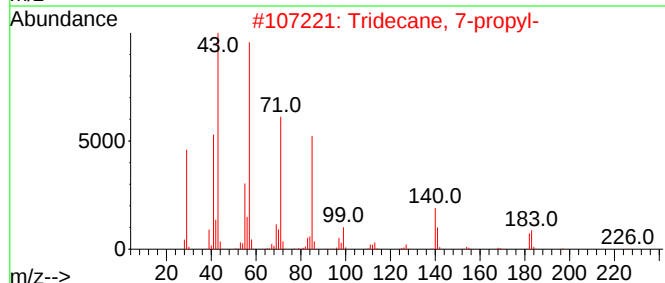
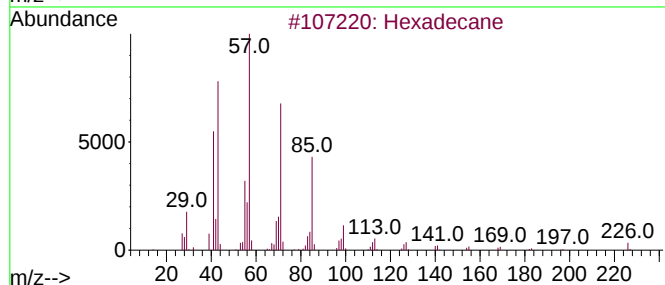
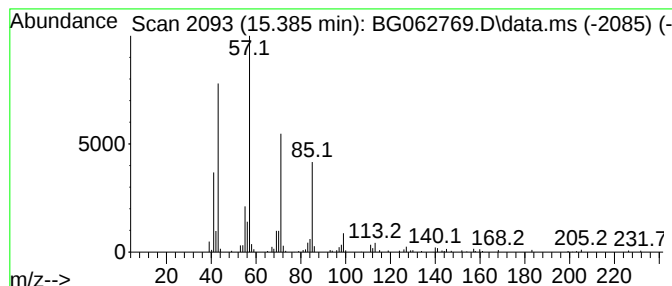
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.385	5.85 ng/ul	259492	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			Tridecane, 7-propyl-	226	C16H34	055045-09-5	90
3			Decane, 3-bromo-	220	C10H21Br	030571-71-2	86
4			Tetradecane	198	C14H30	000629-59-4	86
5			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

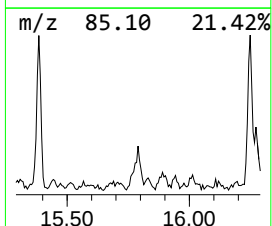
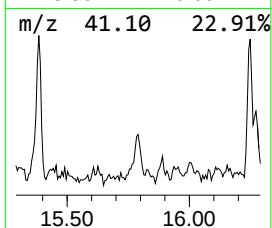
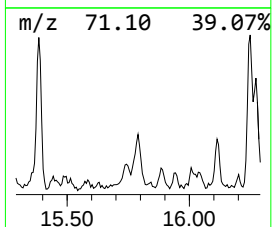
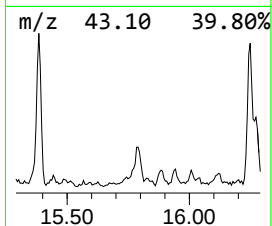
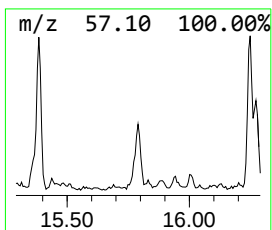
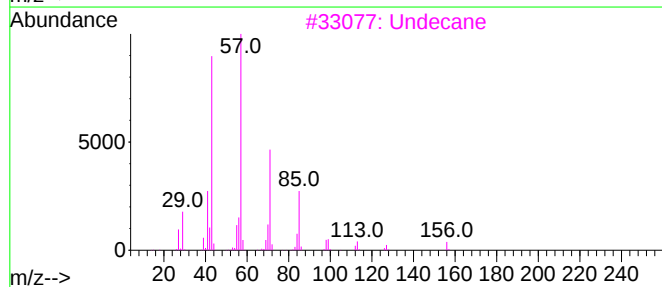
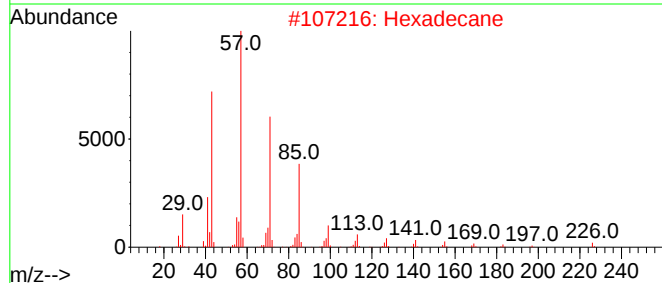
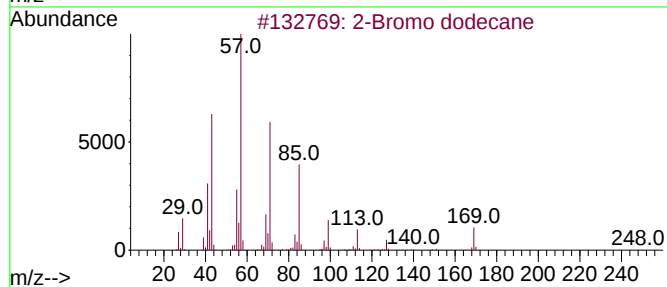
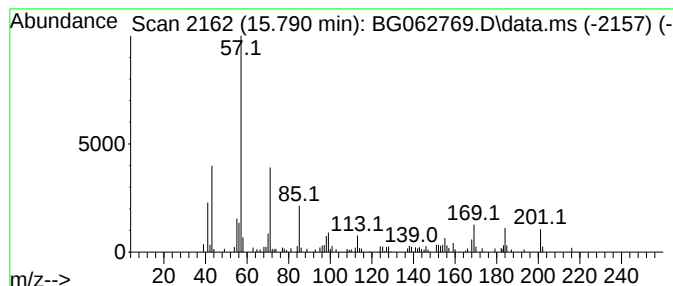
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 2-Bromo dodecane Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.790	2.56 ng/ul	113504	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Bromo dodecane	248	C12H25Br	013187-99-0	59
2			Hexadecane	226	C16H34	000544-76-3	50
3			Undecane	156	C11H24	001120-21-4	47
4			Undecane, 5-ethyl-	184	C13H28	017453-94-0	47
5			Heptadecane	240	C17H36	000629-78-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

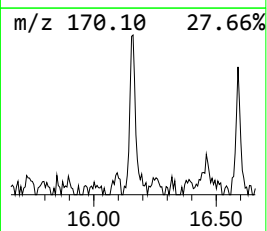
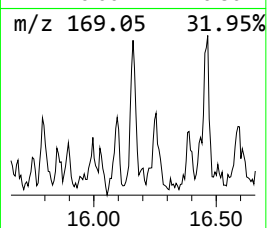
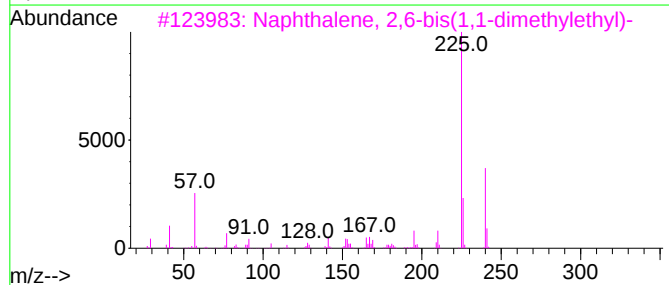
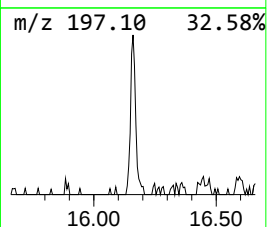
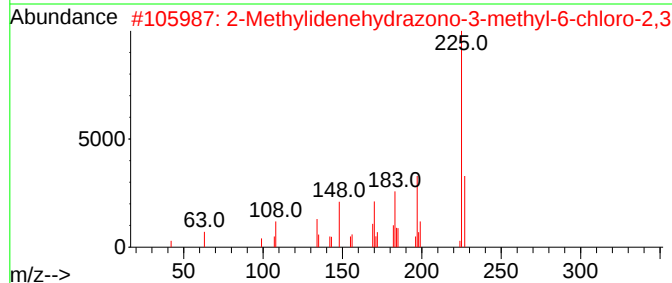
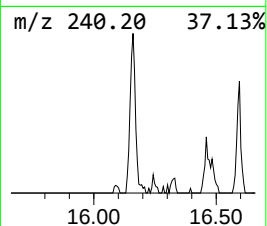
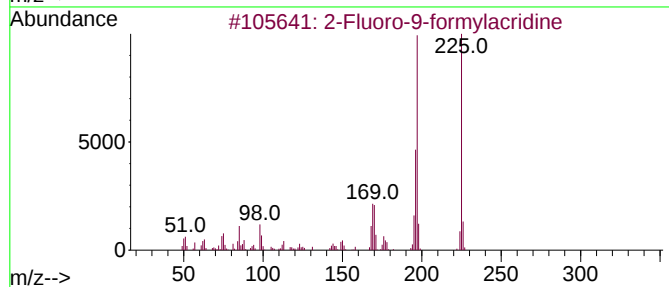
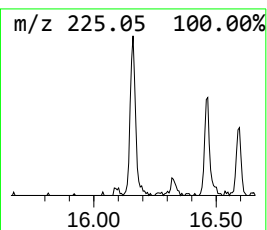
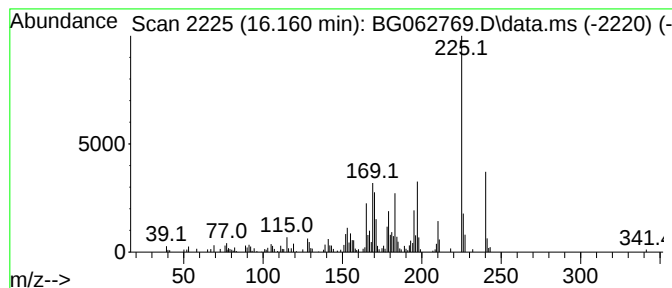
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 unknown-06 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.160	2.15 ng/ul	137254	Phenanthrene-d10	17.276

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Fluoro-9-formylacridine	225	C14H8FNO	024986-45-6	49
2			2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-5	49
3			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	46
4			Phenol, 4-chloro-2,6-bis(1,1-dim...	240	C14H21ClO	004096-72-4	43
5			6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

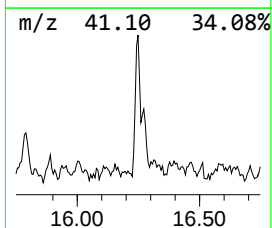
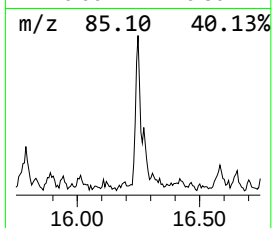
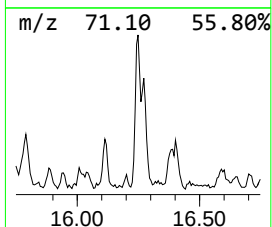
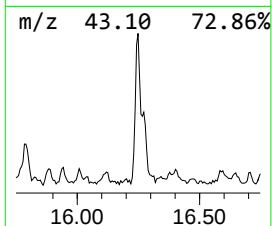
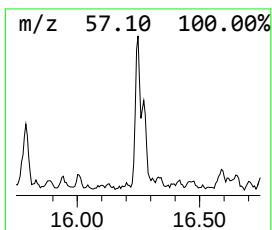
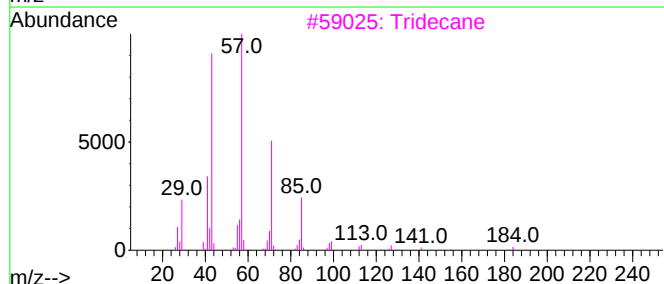
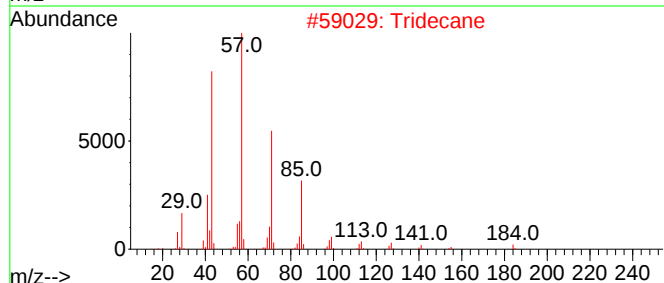
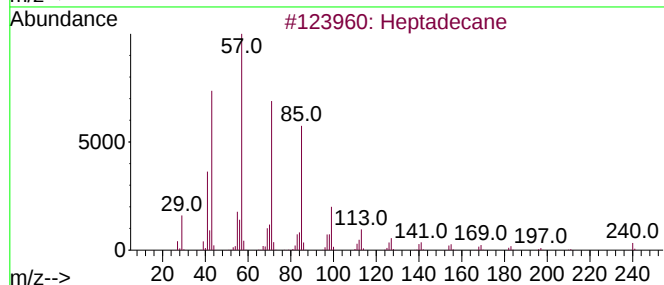
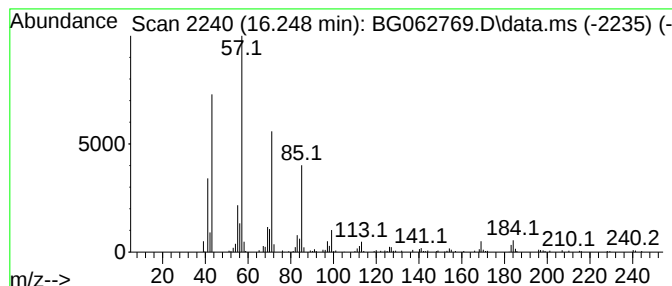
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.248	3.49 ng/ul	223228	Phenanthrene-d10	17.276

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Tridecane	184	C13H28	000629-50-5	90
3		Tridecane	184	C13H28	000629-50-5	87
4		Eicosane	282	C20H42	000112-95-8	86
5		Heneicosane	296	C21H44	000629-94-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

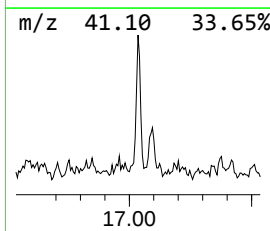
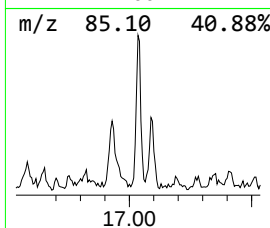
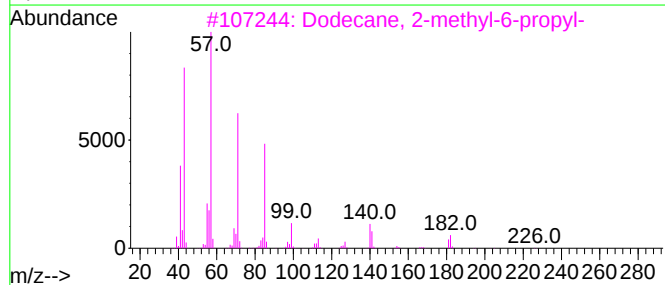
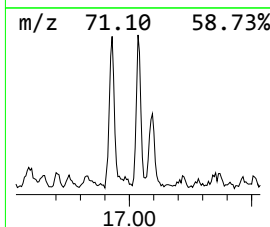
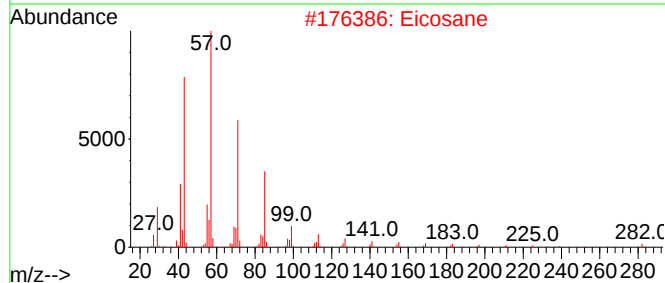
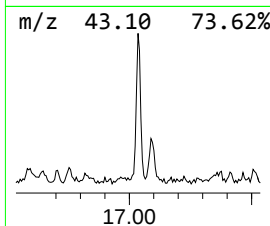
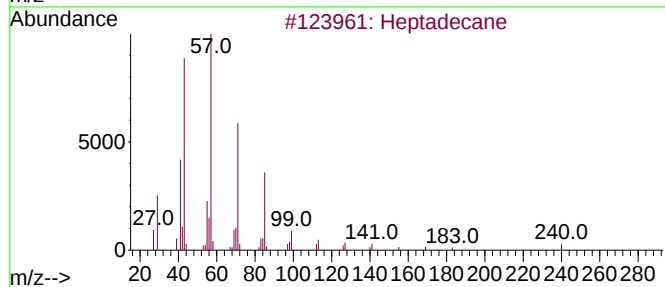
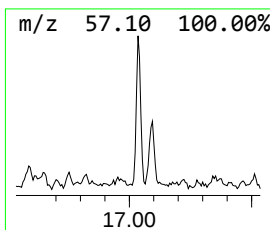
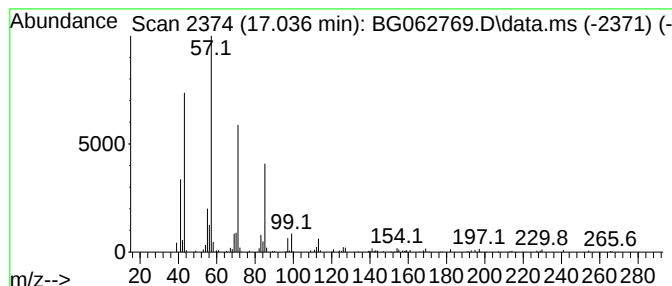
Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

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 47 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.036	3.32 ng/ul	212506	Phenanthrene-d10		17.276	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	83
2	Eicosane		282	C20H42	000112-95-8	83
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	83
4	Heptadecane, 2-methyl-		254	C18H38	001560-89-0	64
5	Heptacosane		380	C27H56	000593-49-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

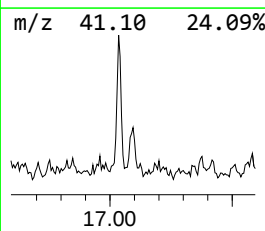
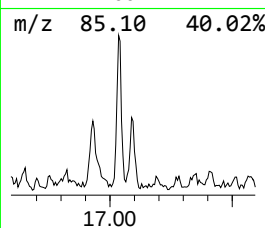
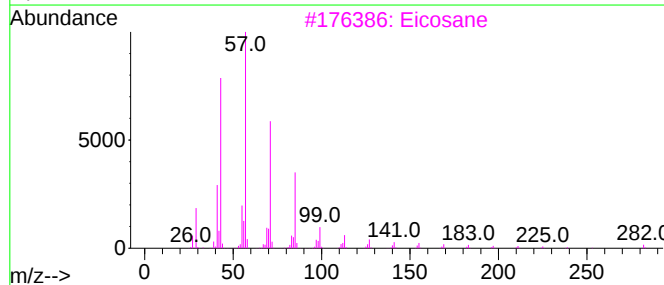
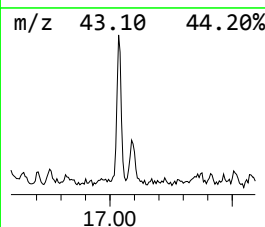
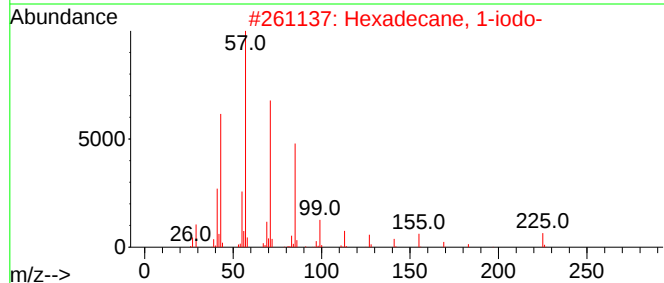
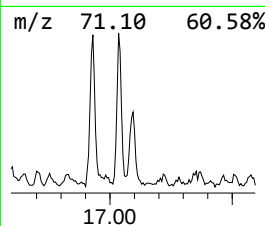
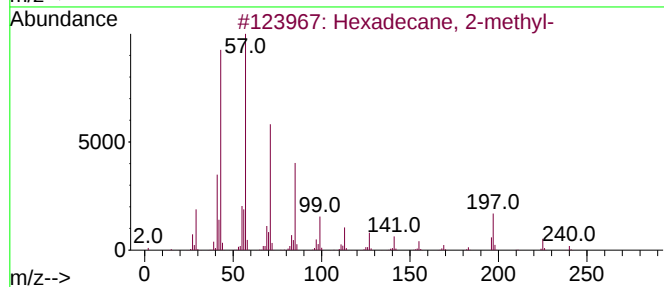
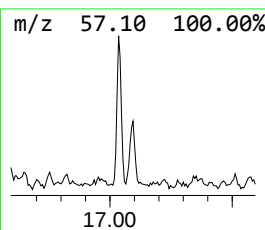
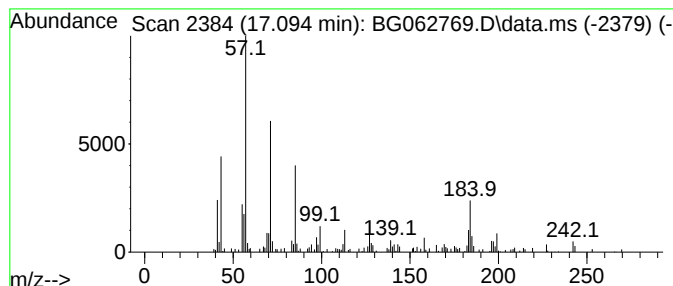
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.094	2.80 ng/ul	178769	Phenanthrene-d10	17.276

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	72
2			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	64
3			Eicosane	282	C20H42	000112-95-8	59
4			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

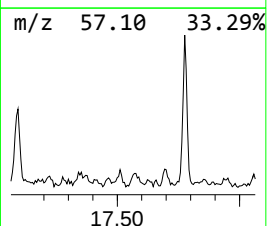
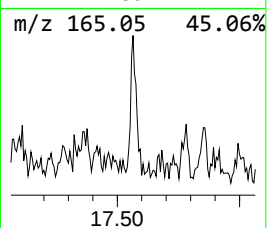
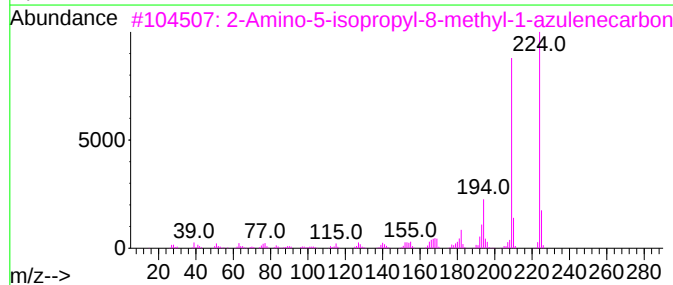
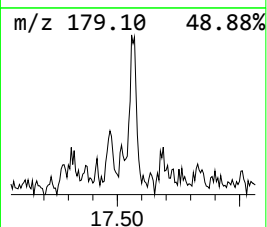
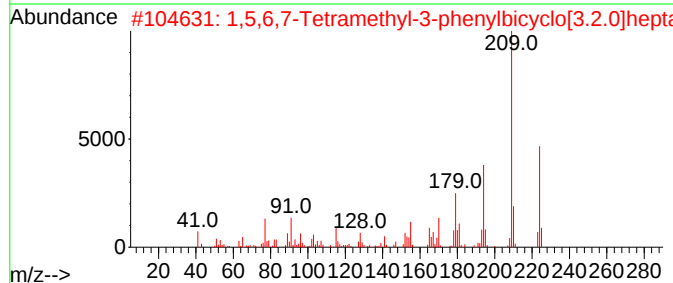
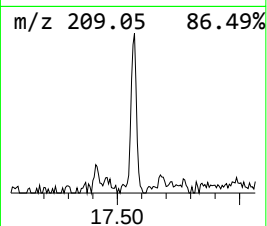
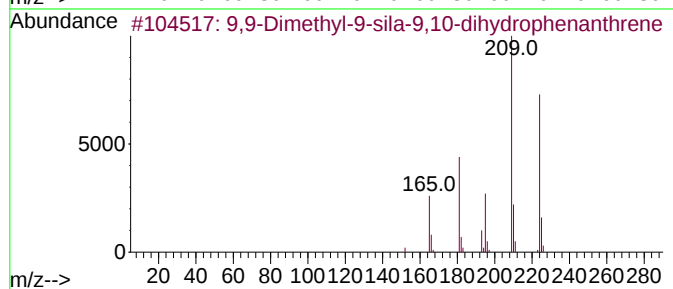
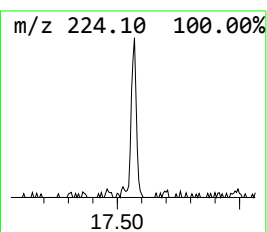
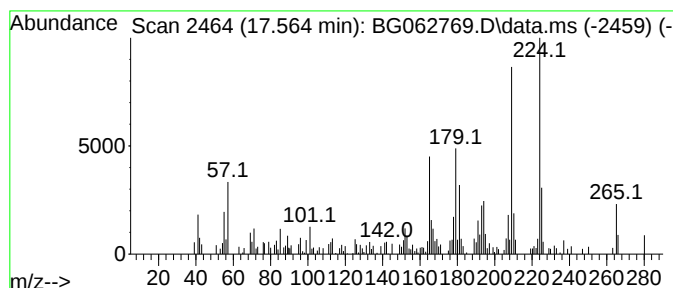
Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

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 49 9,9-Dimethyl-9-sila-9,10-di... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.564	2.08 ng/ul	133084	Phenanthrene-d10	17.276	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si	187328-55-8	62
2	1,5,6,7-Tetramethyl-3-phenylbicy...	224	C17H20	126584-00-7	60
3	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	60
4	Silane, dimethyl(3-ethylphenoxy)...	224	C12H20O2Si	1000347-46-1	58
5	3-(N,N-Dimethylamino)-9-methylca...	224	C15H16N2	094127-15-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

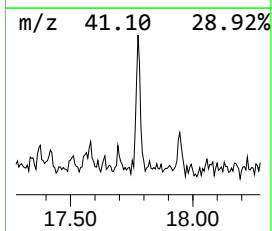
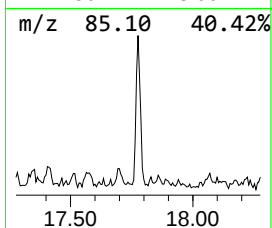
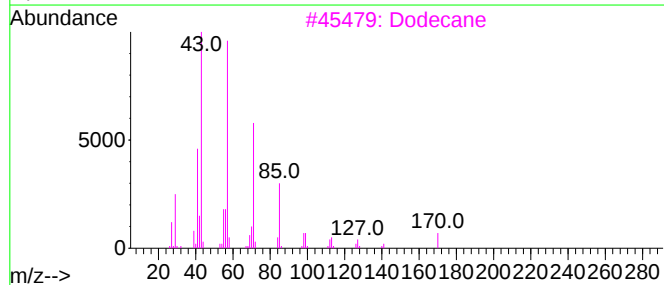
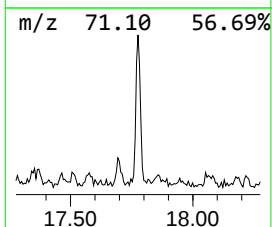
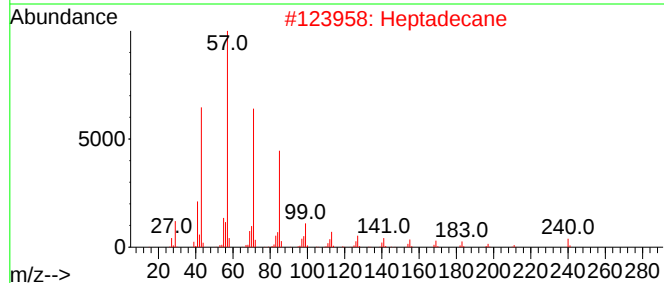
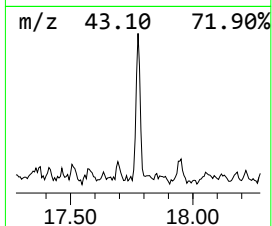
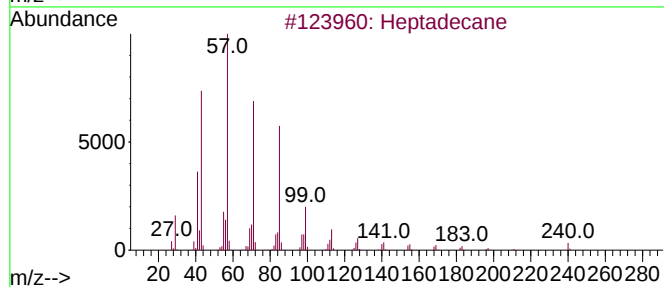
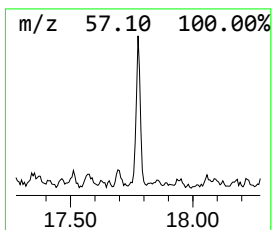
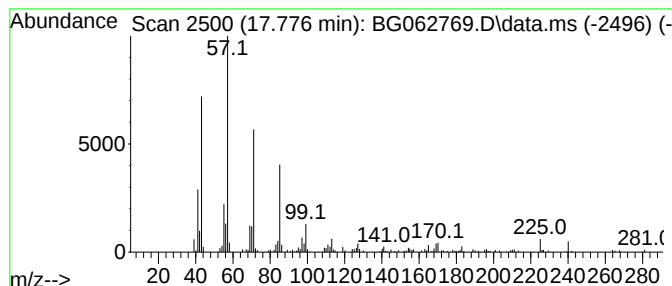
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.776	2.77 ng/ul	176855	Phenanthrene-d10	17.276

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Heptadecane	240	C17H36	000629-78-7	91
3		Dodecane	170	C12H26	000112-40-3	87
4		Octacosane	394	C28H58	000630-02-4	83
5		Heptadecane	240	C17H36	000629-78-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062769.D
Acq On : 12 Sep 2024 19:42
Operator : JU/RC
Sample : P3877-11DL2 30X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVY4DL2

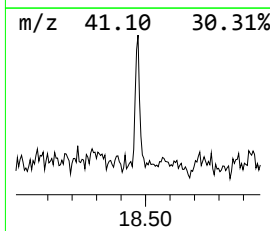
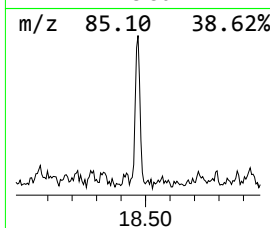
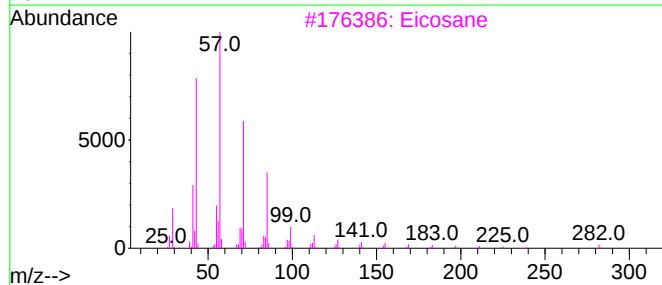
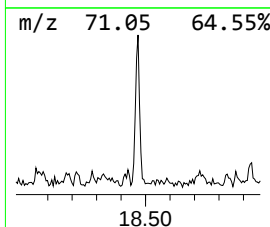
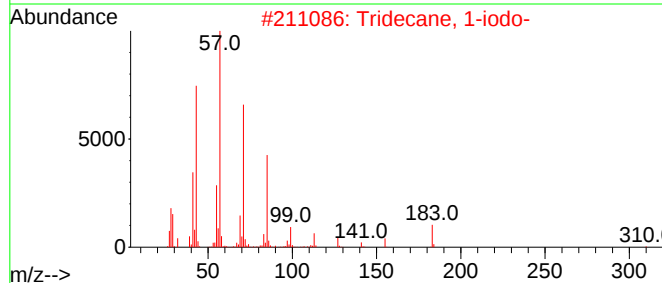
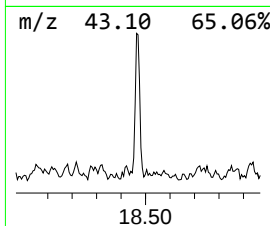
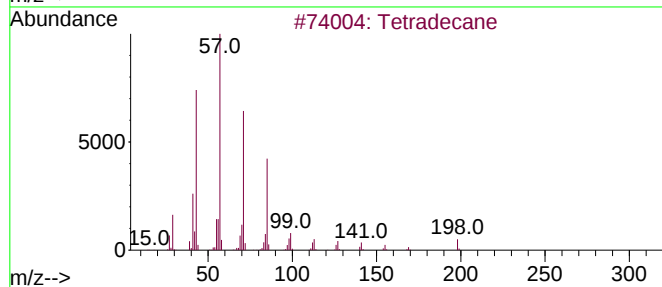
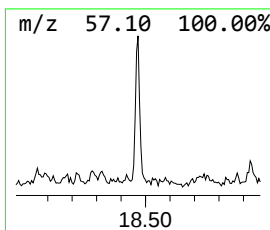
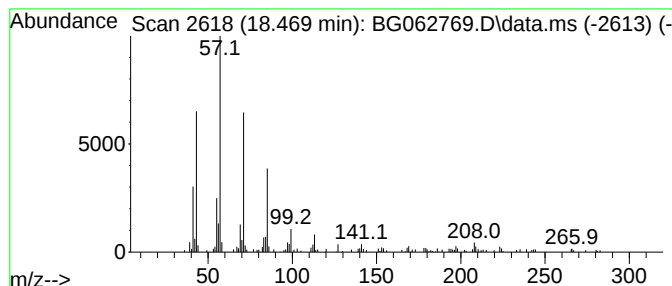
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.469	2.03 ng/ul	129465	Phenanthrene-d10	17.276

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	87
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
3		Eicosane	282	C20H42	000112-95-8	83
4		Tritetracontane	605	C43H88	007098-21-7	80
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
 Data File : BG062769.D
 Acq On : 12 Sep 2024 19:42
 Operator : JU/RC
 Sample : P3877-11DL2 30X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCVY4DL2

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

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.931	12.8	ng/ul	217670	1	7.905	338926	20.0
(DEL) Alkane: C...	6.189	3.0	ng/ul	51700	1	7.905	338926	20.0
(DEL) Alkane: S...	6.466	3.4	ng/ul	58339	1	7.905	338926	20.0
(DEL) Alkane: C...	6.548	4.7	ng/ul	79412	1	7.905	338926	20.0
(DEL) Alkane: S...	6.900	5.3	ng/ul	89210	1	7.905	338926	20.0
(DEL) Alkane: S...	6.959	5.4	ng/ul	90897	1	7.905	338926	20.0
Benzene, 1-ethy...	6.994	4.6	ng/ul	78578	1	7.905	338926	20.0
(DEL) Alkane: S...	7.065	9.5	ng/ul	160392	1	7.905	338926	20.0
Benzene, 1,2,3-...	7.130	3.4	ng/ul	57707	1	7.905	338926	20.0
2,4-Nonadiyne	7.294	3.6	ng/ul	61067	1	7.905	338926	20.0
Butanoic acid, ...	7.417	3.5	ng/ul	59178	1	7.905	338926	20.0
(DEL) Alkane: S...	7.535	35.5	ng/ul	602217	1	7.905	338926	20.0
(DEL) Alkane: S...	7.823	2.8	ng/ul	48024	1	7.905	338926	20.0
1-Hexanol, 2-et...	7.964	7.5	ng/ul	127088	1	7.905	338926	20.0
Benzene, 1,2,4-...	8.023	3.6	ng/ul	61471	1	7.905	338926	20.0
(DEL) Alkane: C...	8.193	3.3	ng/ul	55337	1	7.905	338926	20.0
(DEL) Alkane: S...	8.440	5.9	ng/ul	99445	1	7.905	338926	20.0
(DEL) Alkane: S...	8.499	2.9	ng/ul	49585	1	7.905	338926	20.0
(DEL) Alkane: S...	8.563	10.1	ng/ul	171785	1	7.905	338926	20.0
(DEL) Alkane: S...	8.669	5.6	ng/ul	94572	1	7.905	338926	20.0
Benzeneacetalde...	8.710	3.7	ng/ul	63370	1	7.905	338926	20.0
Benzene, 2-ethy...	8.904	3.7	ng/ul	63080	1	7.905	338926	20.0
Benzene, 4-ethy...	9.004	7.8	ng/ul	131856	1	7.905	338926	20.0
(DEL) Alkane: S...	9.133	29.6	ng/ul	502328	1	7.905	338926	20.0
unknown-01	9.256	6.6	ng/ul	111623	1	7.905	338926	20.0
(DEL) Alkane: C...	9.838	2.3	ng/ul	85385	2	10.696	732835	20.0
(DEL) Alkane: S...	9.979	2.3	ng/ul	83459	2	10.696	732835	20.0
unknown-02	10.813	2.9	ng/ul	106856	2	10.696	732835	20.0
(DEL) Alkane: S...	11.072	6.2	ng/ul	225821	2	10.696	732835	20.0
(DEL) Alkane: S...	11.724	2.7	ng/ul	97739	2	10.696	732835	20.0
Benzene, 1-(2-b...	11.795	5.3	ng/ul	194944	2	10.696	732835	20.0
unknown-03	11.959	3.5	ng/ul	128672	2	10.696	732835	20.0
(DEL) Alkane: S...	12.118	5.2	ng/ul	190334	2	10.696	732835	20.0
unknown-04	12.153	3.0	ng/ul	109823	2	10.696	732835	20.0
(DEL) Alkane: S...	13.358	5.0	ng/ul	223346	3	14.533	886569	20.0
Naphthalene, 2,...	13.710	2.5	ng/ul	110389	3	14.533	886569	20.0
3-Isopropyl-6,1...	14.022	2.3	ng/ul	99709	3	14.533	886569	20.0
Sulfurous acid,...	14.433	10.2	ng/ul	450141	3	14.533	886569	20.0
unknown-05	14.603	2.1	ng/ul	92287	3	14.533	886569	20.0
3,4-Dimethyl(1H...	14.674	6.8	ng/ul	301158	3	14.533	886569	20.0
Naphthalene, 1,...	14.738	2.1	ng/ul	95064	3	14.533	886569	20.0
Methyl 2-diethy...	15.291	7.4	ng/ul	326189	3	14.533	886569	20.0
(DEL) Alkane: S...	15.385	5.8	ng/ul	259492	3	14.533	886569	20.0
2-Bromo dodecane	15.790	2.6	ng/ul	113504	3	14.533	886569	20.0
unknown-06	16.160	2.1	ng/ul	137254	4	17.276	1278440	20.0
(DEL) Alkane: S...	16.248	3.5	ng/ul	223228	4	17.276	1278440	20.0
(DEL) Alkane: S...	17.036	3.3	ng/ul	212506	4	17.276	1278440	20.0
(DEL) Alkane: S...	17.094	2.8	ng/ul	178769	4	17.276	1278440	20.0
9,9-Dimethyl-9-...	17.564	2.1	ng/ul	133084	4	17.276	1278440	20.0
(DEL) Alkane: S...	17.776	2.8	ng/ul	176855	4	17.276	1278440	20.0
(DEL) Alkane: S...	18.469	2.0	ng/ul	129465	4	17.276	1278440	20.0