

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062908.D
 Acq On : 18 Sep 2024 14:29
 Operator : JU/RC
 Sample : P3878-12ME
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCVZ0ME

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-BG091724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062908.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	107	111	120	rBV2	63801	119547	2.69%	0.474%
2	4.980	316	322	332	rVB2	19173	36170	0.81%	0.143%
3	5.891	470	477	484	rVV2	67131	110315	2.48%	0.437%
4	6.173	512	525	531	rBV2	71399	144703	3.26%	0.573%
5	6.361	552	557	563	rVB6	27812	52655	1.19%	0.209%
6	6.425	564	568	573	rVB2	30058	39135	0.88%	0.155%
7	6.502	573	581	585	rBV6	22334	55869	1.26%	0.221%
8	6.766	618	626	631	rVB4	15801	37638	0.85%	0.149%
9	6.854	637	641	646	rVV4	44341	78028	1.76%	0.309%
10	6.913	646	651	654	rVV3	53491	79075	1.78%	0.313%
11	6.954	654	658	662	rVV2	40158	68876	1.55%	0.273%
12	7.025	662	670	677	rVV2	198698	387322	8.72%	1.534%
13	7.083	677	680	685	rVB2	29683	45833	1.03%	0.182%
14	7.189	692	698	703	rBV	100552	172935	3.89%	0.685%
15	7.248	703	708	711	rVV3	25589	42166	0.95%	0.167%
16	7.289	711	715	718	rVB2	31073	44645	1.00%	0.177%
17	7.389	726	732	743	rVB3	131825	274613	6.18%	1.088%
18	7.489	743	749	757	rBV2	267239	495105	11.14%	1.961%
19	7.853	804	811	817	rBV2	196900	370673	8.34%	1.468%
20	7.918	817	822	828	rVV2	127225	204565	4.60%	0.810%
21	7.976	828	832	839	rVB5	29120	55450	1.25%	0.220%
22	8.147	859	861	868	rVB6	26575	40242	0.91%	0.159%
23	8.282	879	884	888	rBV3	39874	69688	1.57%	0.276%
24	8.394	898	903	909	rVB3	44405	84810	1.91%	0.336%
25	8.452	909	913	916	rVB3	30042	40340	0.91%	0.160%
26	8.499	916	921	922	rBV2	57368	84161	1.89%	0.333%
27	8.558	927	931	938	rVV2	174350	290931	6.55%	1.152%
28	8.623	938	942	945	rVV2	48046	68129	1.53%	0.270%
29	8.658	945	948	956	rVB2	37601	67490	1.52%	0.267%
30	8.811	968	974	978	rBV	28617	48231	1.09%	0.191%
31	8.858	978	982	989	rVB	35361	57601	1.30%	0.228%
32	8.963	990	1000	1002	rVV5	48540	99029	2.23%	0.392%
33	9.005	1003	1007	1013	rVV	137169	242887	5.47%	0.962%
34	9.087	1016	1021	1026	rVB	243557	370695	8.34%	1.468%
35	9.210	1032	1042	1048	rVB9	36299	104863	2.36%	0.415%
36	9.645	1113	1116	1122	rVB3	27390	40382	0.91%	0.160%

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 ClientSampleId :
 DCVZ0ME

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-BG091724.MA.M
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37	9.727	1122	1130	1136	rBV4	140574	285976	6.44%	1.133%
38	9.792	1136	1141	1145	rBV5	32248	66617	1.50%	0.264%
39	9.933	1162	1165	1169	rVB4	37742	51474	1.16%	0.204%
40	9.998	1171	1176	1181	rBV8	24816	43317	0.97%	0.172%
41	10.268	1216	1222	1230	rVB2	183900	335344	7.55%	1.328%
42	10.338	1230	1234	1243	rBV5	31788	58451	1.32%	0.232%
43	10.638	1276	1285	1292	rBV2	336916	772289	17.38%	3.059%
44	10.779	1301	1309	1319	rBV3	183994	420516	9.46%	1.666%
45	10.926	1329	1334	1343	rVB7	29032	70588	1.59%	0.280%
46	11.020	1345	1350	1358	rBV	131233	221975	5.00%	0.879%
47	11.149	1365	1372	1380	rBV3	53910	111453	2.51%	0.441%
48	11.566	1435	1443	1451	rBV3	22516	65614	1.48%	0.260%
49	11.672	1456	1461	1466	rVV2	57749	99522	2.24%	0.394%
50	11.743	1466	1473	1482	rVB8	32906	86842	1.95%	0.344%
51	11.913	1497	1502	1511	rVB2	69750	114182	2.57%	0.452%
52	12.072	1522	1529	1532	rBV2	104399	174875	3.94%	0.693%
53	12.107	1532	1535	1541	rVB2	80039	111294	2.50%	0.441%
54	12.313	1564	1570	1577	rVB3	137215	245470	5.52%	0.972%
55	12.477	1593	1598	1602	rBV6	31981	57910	1.30%	0.229%
56	12.524	1602	1606	1615	rVV7	33582	69125	1.56%	0.274%
57	12.888	1663	1668	1675	rVB8	24375	42207	0.95%	0.167%
58	13.029	1688	1692	1697	rBV2	29508	44636	1.00%	0.177%
59	13.317	1736	1741	1748	rVB	135403	204750	4.61%	0.811%
60	13.670	1793	1801	1805	rBV5	46243	103606	2.33%	0.410%
61	13.746	1810	1814	1819	rBV4	29633	45181	1.02%	0.179%
62	13.811	1821	1825	1830	rVV6	47083	115283	2.59%	0.457%
63	13.893	1834	1839	1845	rVB	352781	529362	11.91%	2.097%
64	13.975	1851	1853	1857	rVB3	54454	61582	1.39%	0.244%
65	14.181	1882	1888	1898	rVB	364094	578479	13.02%	2.291%
66	14.392	1920	1924	1929	rVB	189244	238515	5.37%	0.945%
67	14.486	1935	1940	1948	rVV	477050	756524	17.03%	2.997%
68	14.569	1950	1954	1960	rVB7	51660	89512	2.01%	0.355%
69	14.633	1960	1965	1968	rBV2	45507	62391	1.40%	0.247%
70	14.692	1970	1975	1985	rBV3	162279	298825	6.73%	1.184%
71	15.021	2027	2031	2037	rBV8	29808	58565	1.32%	0.232%
72	15.115	2042	2047	2058	rVB2	260024	469192	10.56%	1.858%
73	15.250	2067	2070	2077	rBV7	31558	81084	1.82%	0.321%
74	15.344	2083	2086	2092	rVB	135847	184885	4.16%	0.732%
75	15.485	2104	2110	2115	rVB	475264	733187	16.50%	2.904%
76	15.597	2124	2129	2137	rVB2	146498	244405	5.50%	0.968%

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 ClientSampleId :
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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

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 Title : SVOA CALIBRATION

77	15.850	2167	2172	2177	rVB2	96569	151949	3.42%	0.602%
78	16.208	2229	2233	2236	rBV	165982	217593	4.90%	0.862%
79	16.549	2287	2291	2294	rBV6	27564	38774	0.87%	0.154%
80	16.895	2343	2350	2362	rBV	2822804	4443119	100.00%	17.599%
81	17.001	2365	2368	2372	rVV2	142872	180099	4.05%	0.713%
82	17.054	2374	2377	2382	rVB2	74370	101786	2.29%	0.403%
83	17.242	2403	2409	2414	rBV	613101	917196	20.64%	3.633%
84	17.336	2420	2425	2430	rVB	559406	792408	17.83%	3.139%
85	17.741	2490	2494	2499	rVB2	118122	150304	3.38%	0.595%
86	17.824	2504	2508	2510	rVV5	31581	43134	0.97%	0.171%
87	17.853	2511	2513	2518	rVV5	48296	60797	1.37%	0.241%
88	17.912	2520	2523	2527	rVB2	54079	61437	1.38%	0.243%
89	18.435	2608	2612	2616	rBV	96412	116780	2.63%	0.463%
90	19.069	2716	2720	2726	rVB3	83338	163456	3.68%	0.647%
91	19.633	2810	2816	2822	rVB	736636	1037950	23.36%	4.111%
92	20.180	2906	2909	2914	rVB3	49167	62441	1.41%	0.247%
93	20.673	2990	2993	2997	rBV4	38206	38483	0.87%	0.152%
94	21.161	3071	3076	3082	rVB3	210285	313236	7.05%	1.241%
95	21.490	3126	3132	3138	rVB	631188	997984	22.46%	3.953%
96	21.684	3163	3165	3169	rVB3	33983	38464	0.87%	0.152%
97	22.142	3239	3243	3249	rVB4	68222	110892	2.50%	0.439%
98	22.266	3259	3264	3270	rVB7	46252	94722	2.13%	0.375%
99	24.316	3605	3613	3623	rVB	382045	977289	22.00%	3.871%
100	24.528	3640	3649	3661	rVB2	415626	1105844	24.89%	4.380%

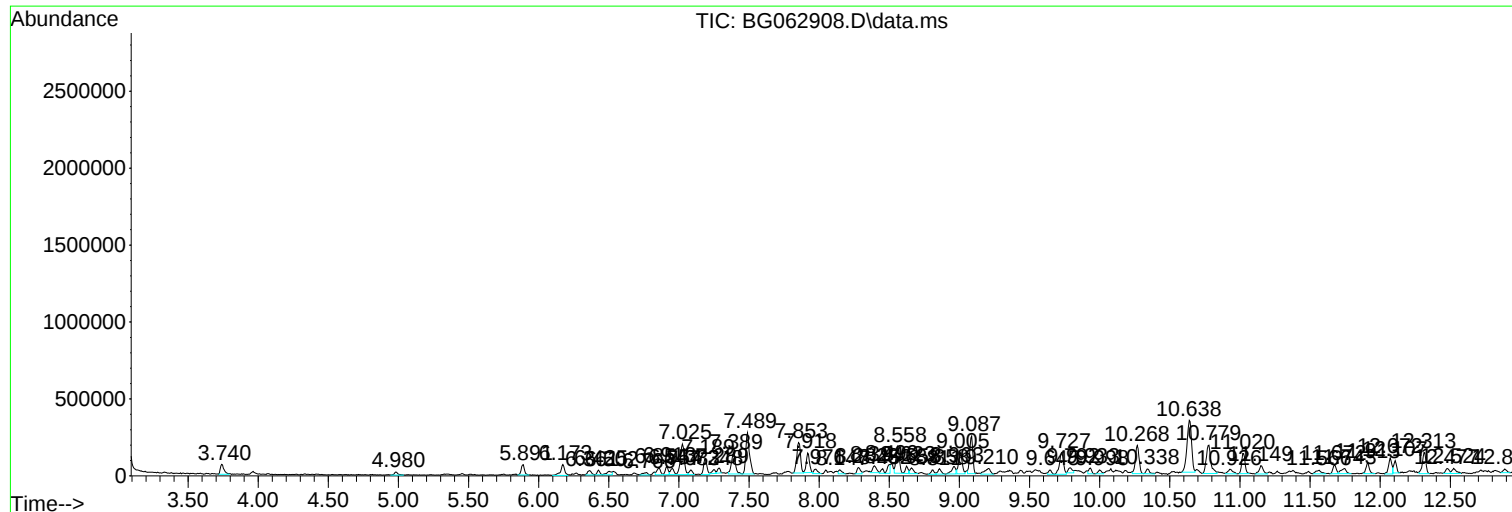
Sum of corrected areas: 25245939

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ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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



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Misc :
ALS Vial : 8 Sample Multiplier: 1

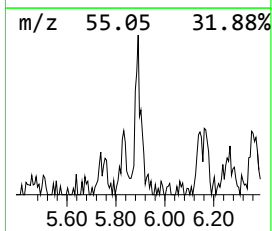
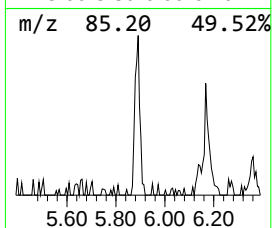
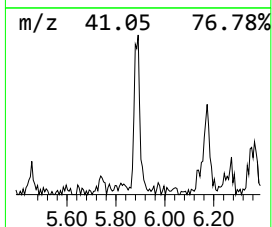
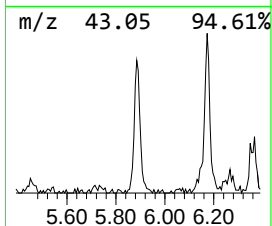
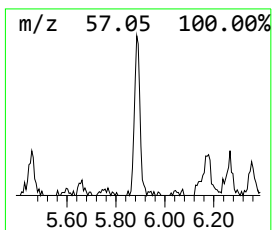
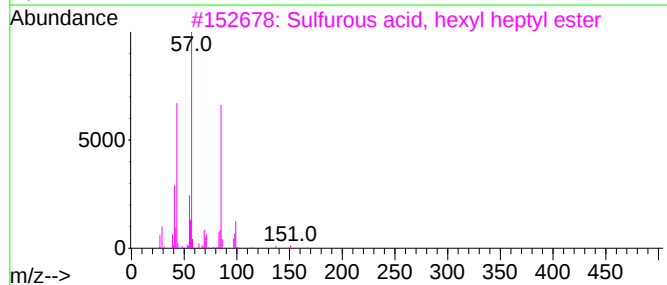
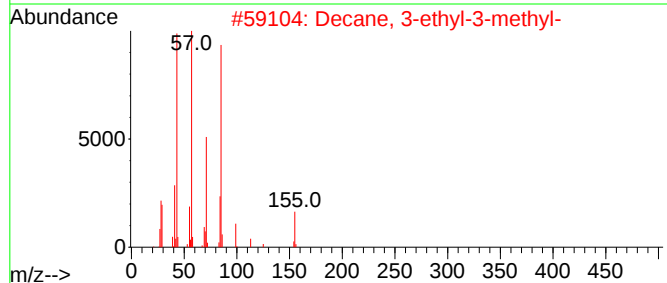
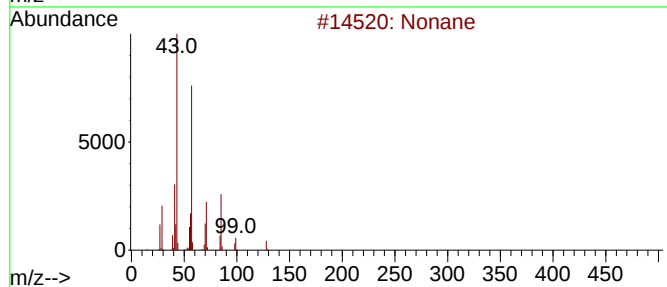
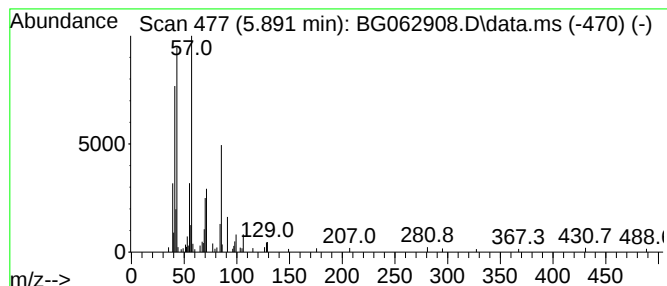
Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.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 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.891	5.95 ng/ul	110315	1,4-Dichlorobenzene-d4			7.859
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	52
2	Decane, 3-ethyl-3-methyl-		184	C13H28	017312-66-2	50
3	Sulfurous acid, hexyl heptyl ester		264	C13H28O3S	1000309-12-9	47
4	Decane, 1-iodo-		268	C10H21I	002050-77-3	47
5	2,4-Dimethyldodecane		198	C14H30	006117-99-3	47



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Instrument :
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ClientSampleId :
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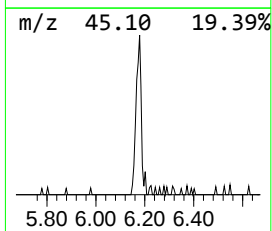
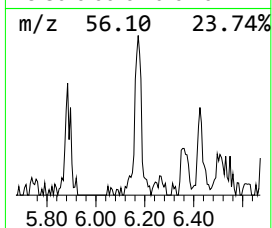
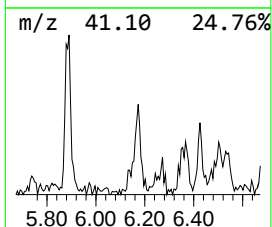
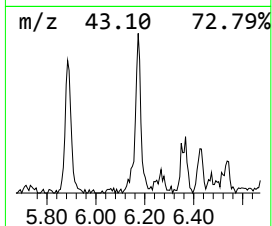
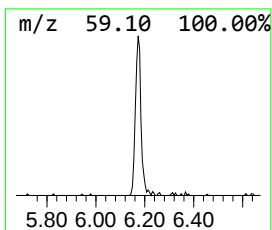
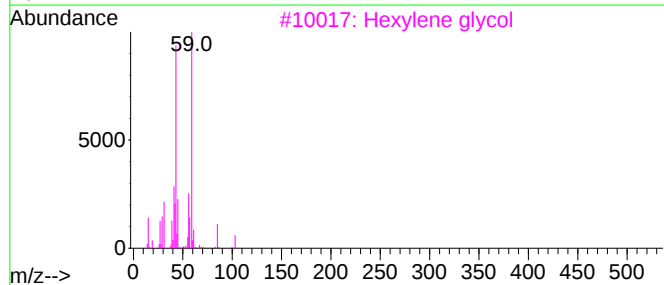
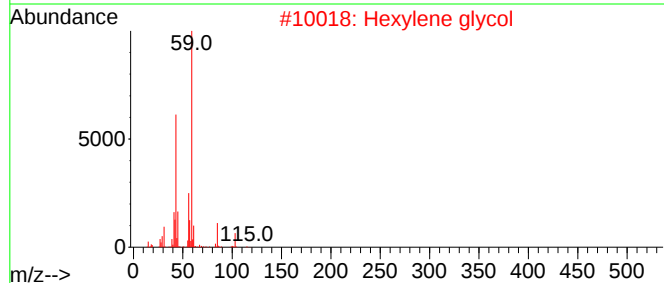
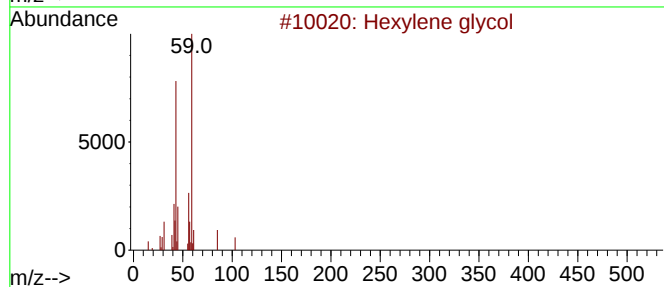
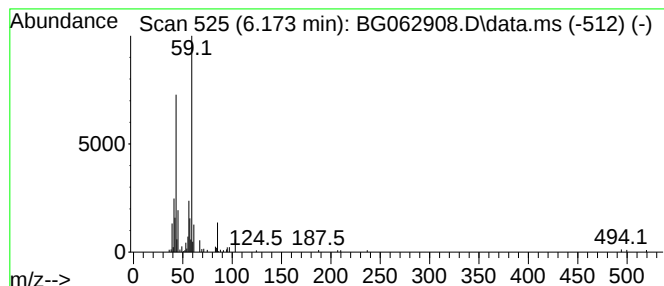
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.173	7.81 ng/ul	144703	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	90
3			Hexylene glycol	118	C6H14O2	000107-41-5	83
4			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	83
5			Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	64



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ClientSampleId :
DCVZ0ME

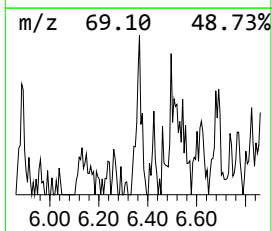
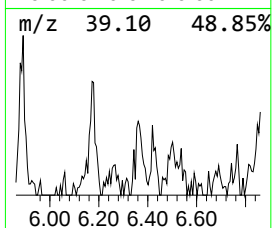
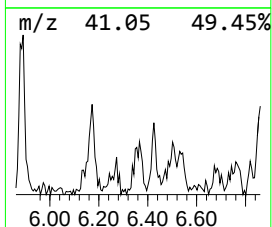
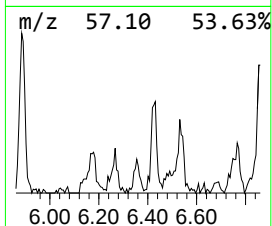
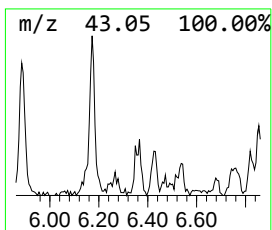
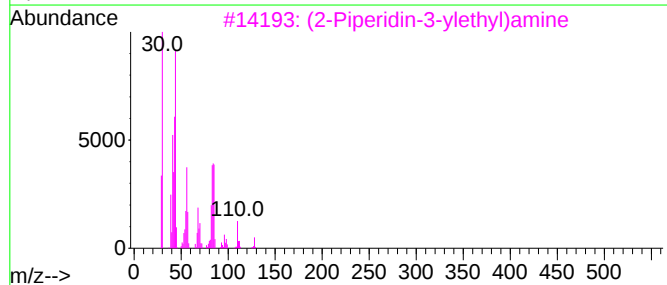
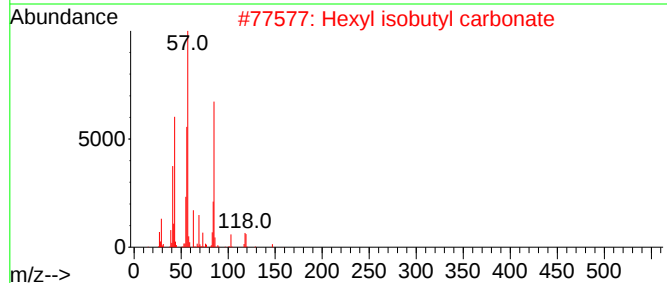
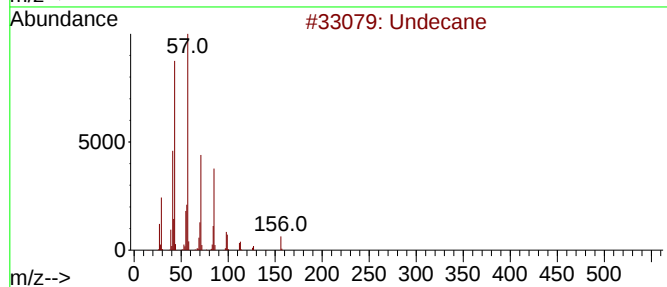
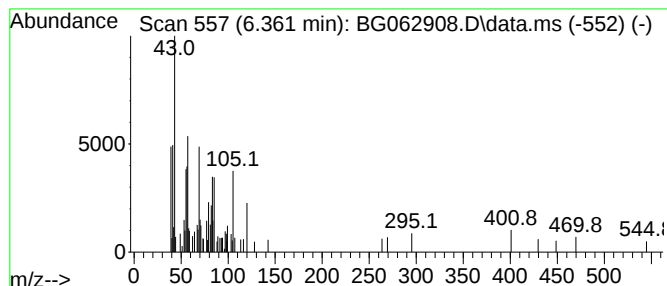
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.361	2.84 ng/ul	52655	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	38
2			Hexyl isobutyl carbonate	202	C11H22O3	959067-98-2	38
3			(2-Piperidin-3-ylethyl)amine	128	C7H16N2	089850-94-2	38
4			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	38
5			Oxalic acid, isobutyl hexyl ester	230	C12H22O4	1000309-37-1	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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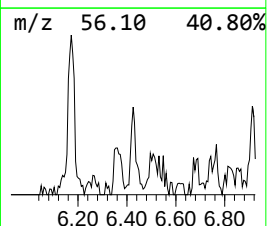
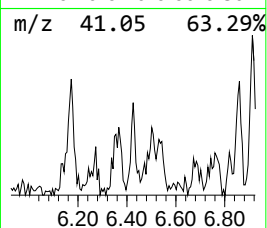
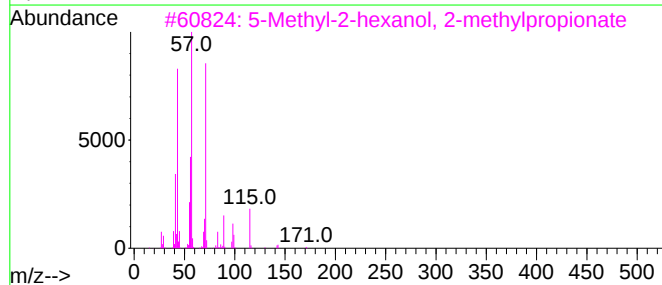
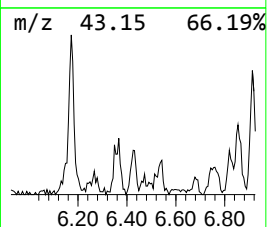
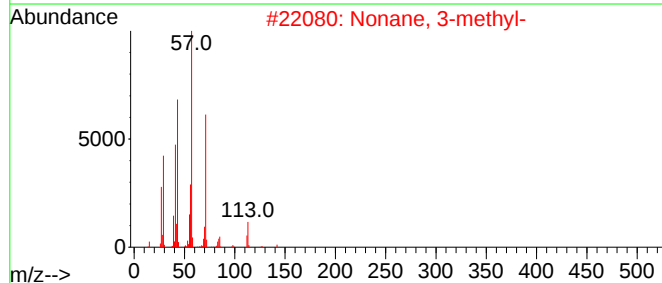
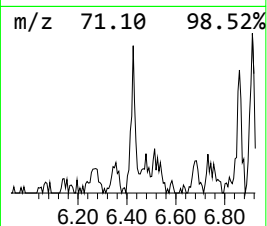
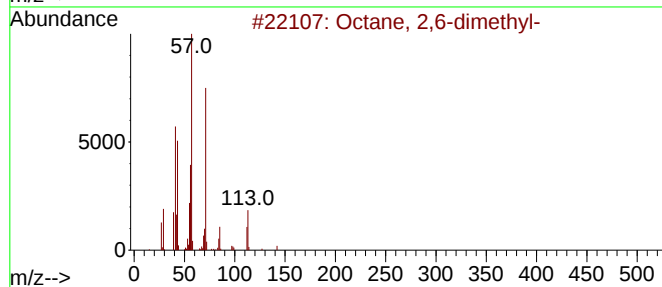
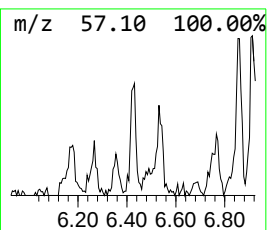
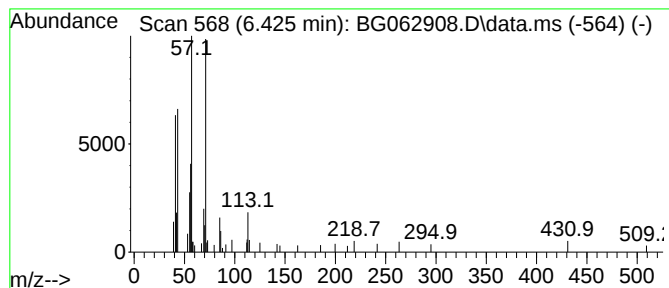
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.425	2.11 ng/ul	39135	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	81
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	64
3			5-Methyl-2-hexanol, 2-methylprop...	186	C11H22O2	1000447-34-4	53
4			Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	53
5			6-Methyl-2-heptanol, 2-methylpro...	200	C12H24O2	1000447-36-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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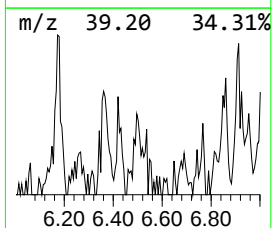
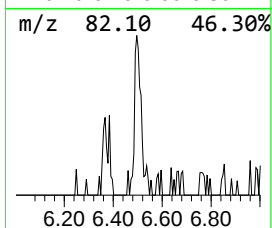
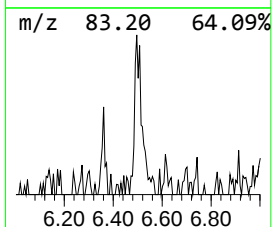
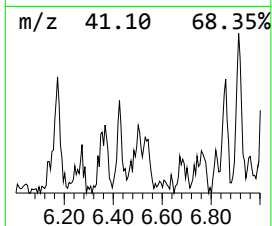
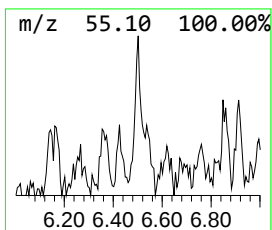
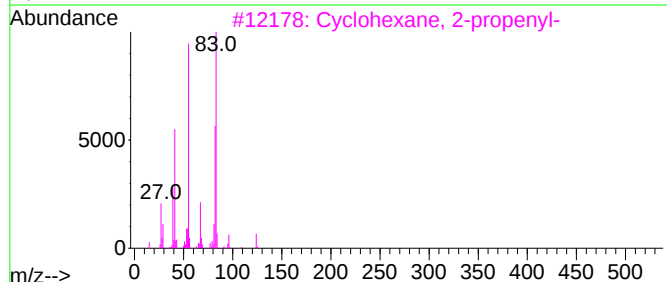
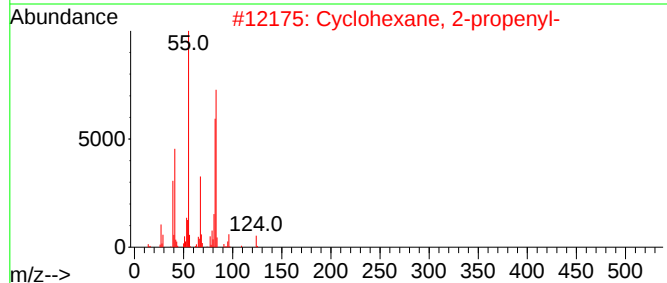
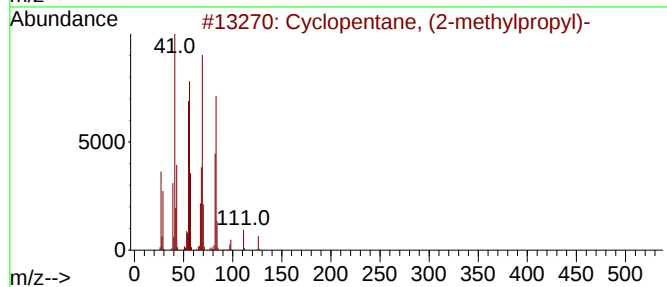
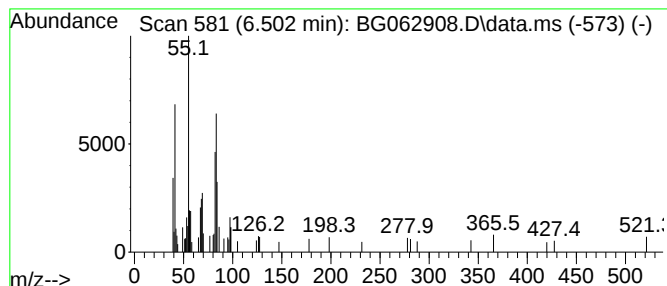
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic6.502 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.502	3.01 ng/ul	55869	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, (2-methylpropyl)-	126	C9H18	003788-32-7	47
2			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	43
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	43
4			Piperidine, 4-(4-(2-chlorophenyl...	261	C14H16ClN3	902836-42-4	43
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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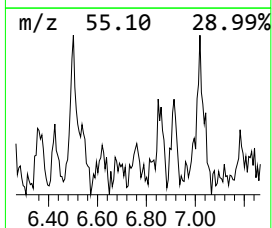
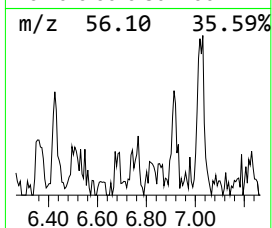
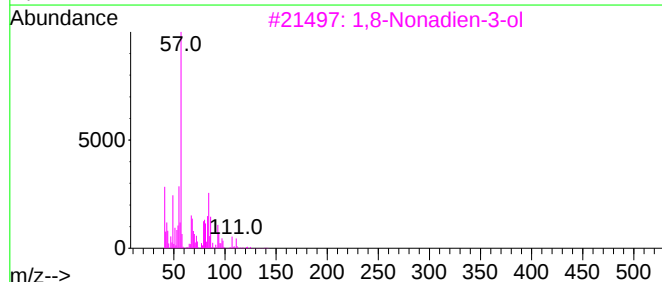
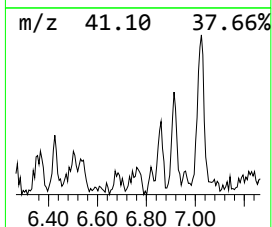
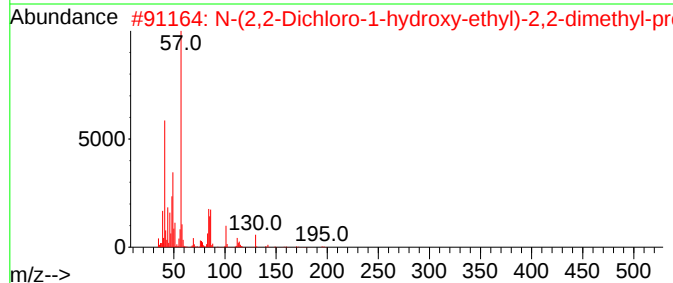
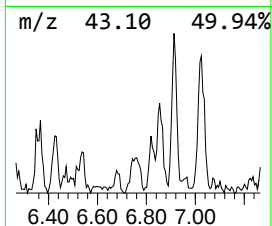
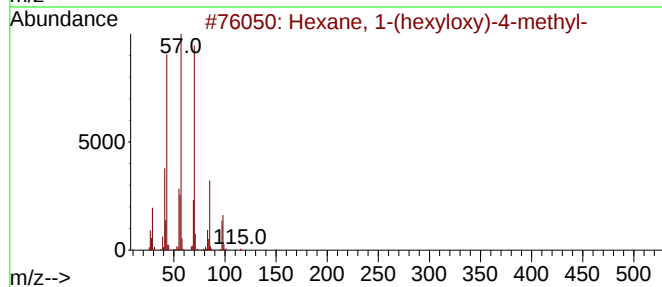
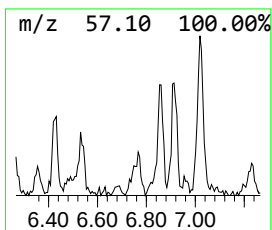
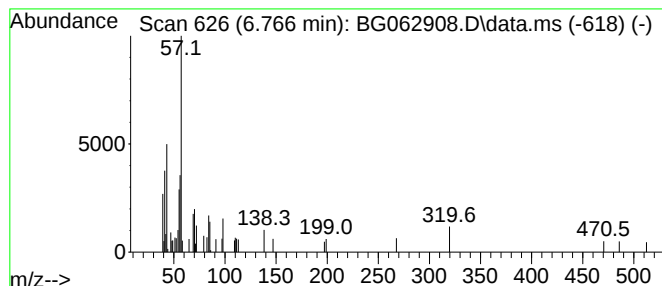
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.766	2.03 ng/ul	37638	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1-(hexyloxy)-4-methyl-	200	C13H28O	074421-20-8	9
2			N-(2,2-Dichloro-1-hydroxy-ethyl)...	213	C7H13Cl2NO2	058956-78-8	9
3			1,8-Nonadien-3-ol	140	C9H16O	159010-02-3	9
4			Propanamide, 2,2-dimethyl-N-(3-n...	222	C11H14N2O3	032597-30-1	9
5			Diethylmalonic acid, monochlorid...	318	C17H31ClO3	1000369-41-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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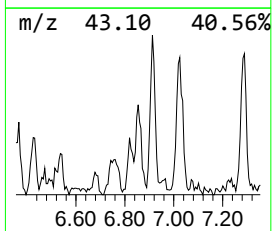
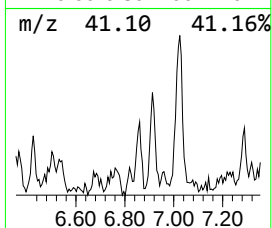
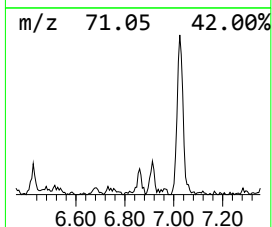
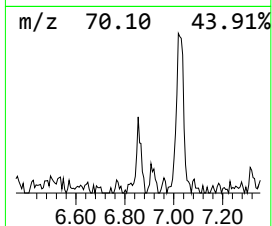
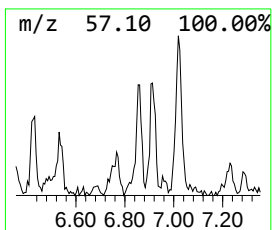
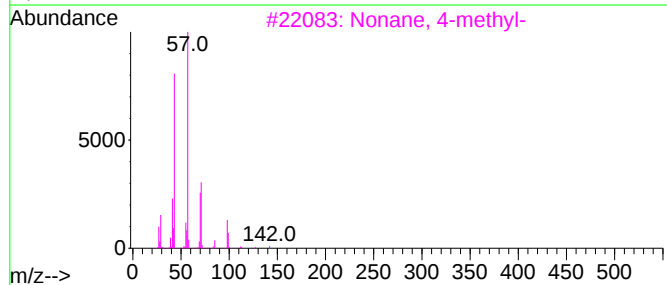
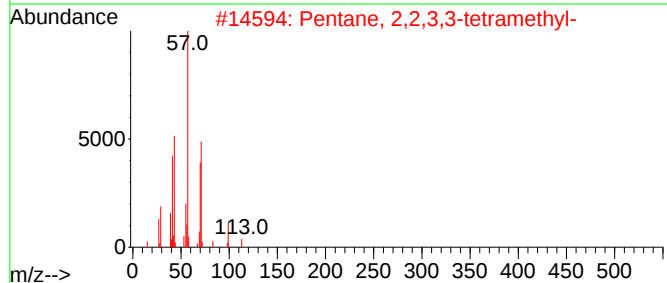
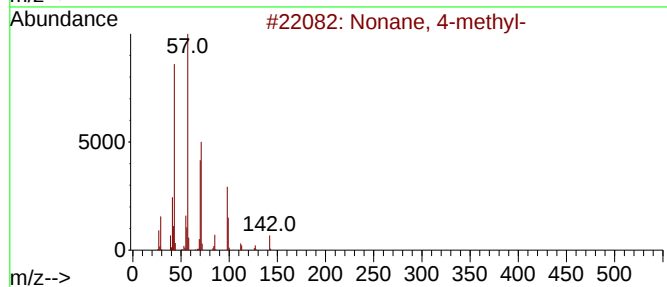
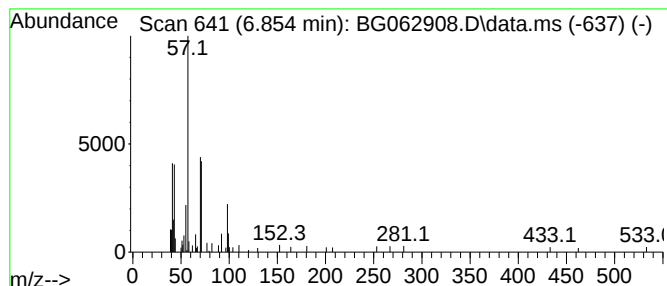
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.854	4.21 ng/ul	78028	1,4-Dichlorobenzene-d4	7.859

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 4-methyl-	142	C10H22	017301-94-9	50
2		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	45
3		Nonane, 4-methyl-	142	C10H22	017301-94-9	45
4		Nonane, 4-methyl-	142	C10H22	017301-94-9	40
5		1-Decanol, 2-methyl-	172	C11H24O	018675-24-6	39



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

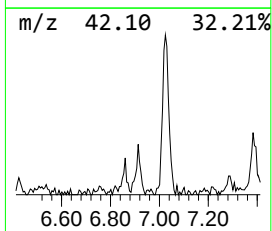
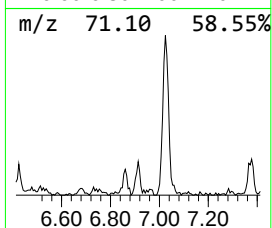
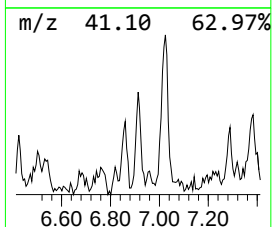
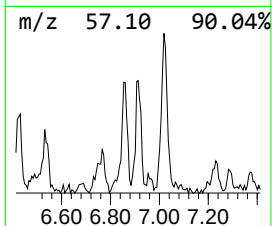
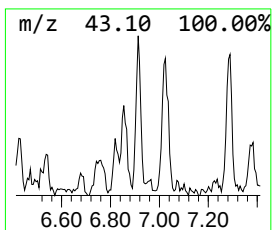
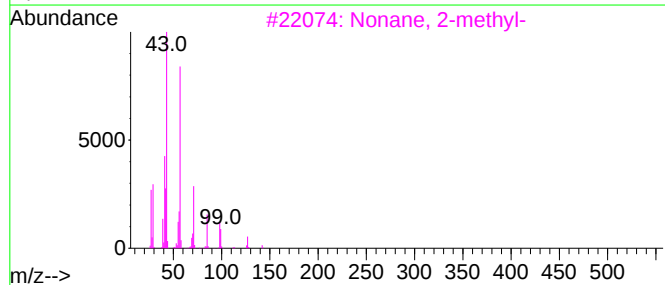
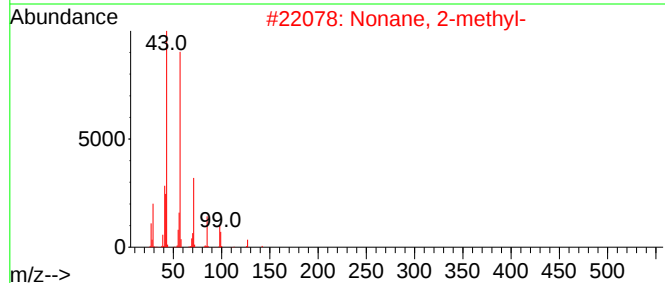
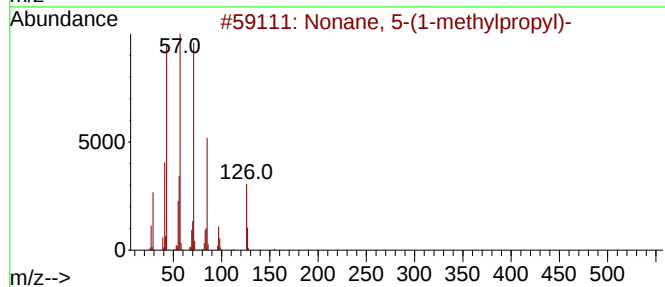
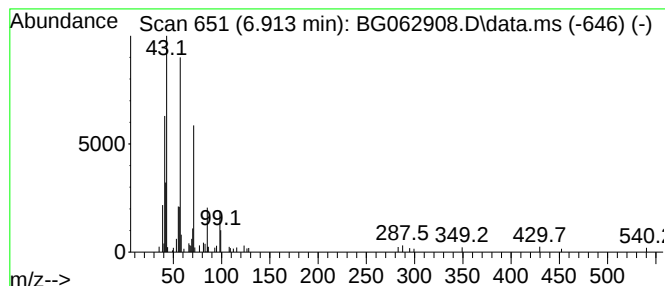
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.913	4.27 ng/ul	79075	1,4-Dichlorobenzene-d4			7.859
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 5-(1-methylpropyl)-		184	C13H28	062185-54-0	59
2	Nonane, 2-methyl-		142	C10H22	000871-83-0	59
3	Nonane, 2-methyl-		142	C10H22	000871-83-0	59
4	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	59
5	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
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ClientSampleId :
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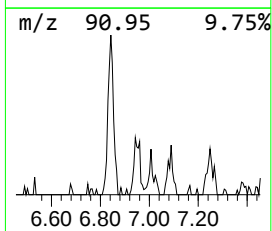
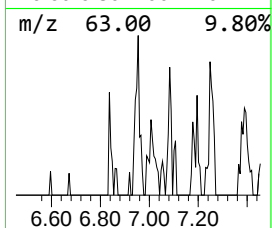
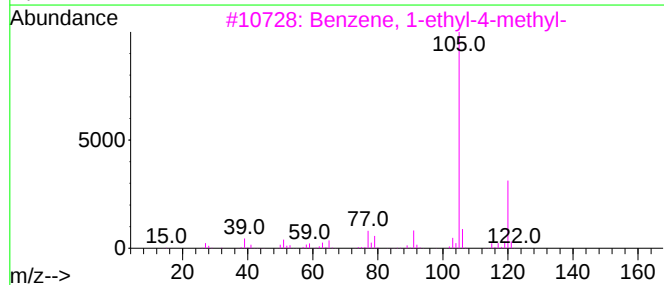
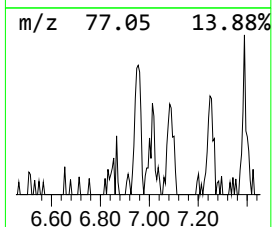
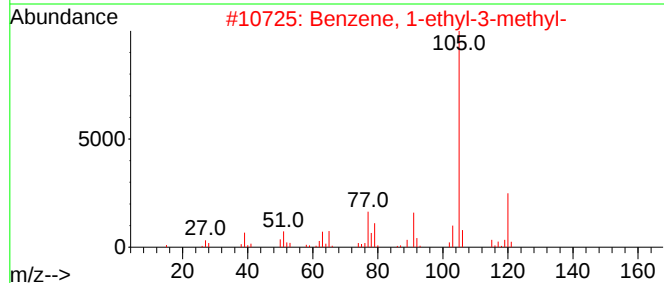
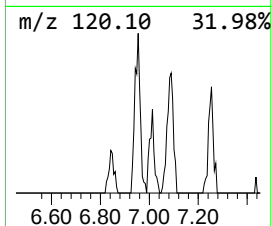
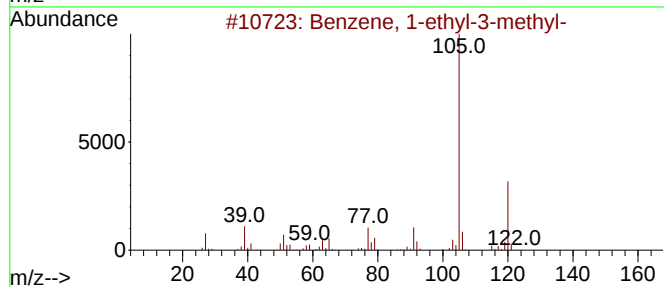
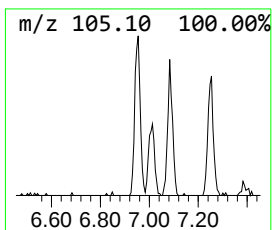
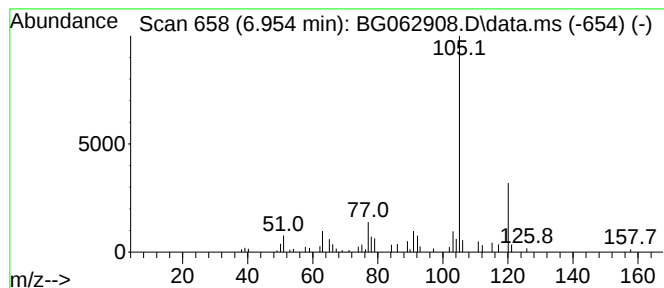
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.954	3.72 ng/ul	68876	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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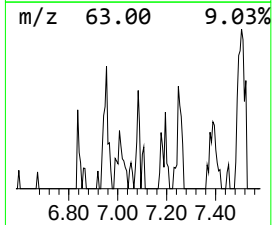
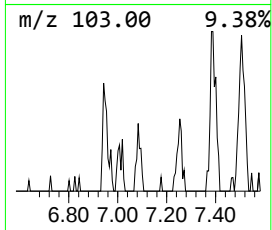
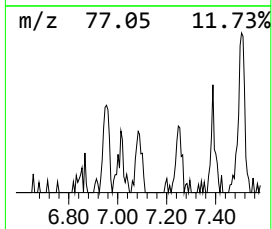
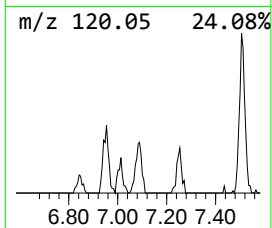
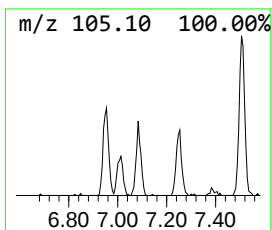
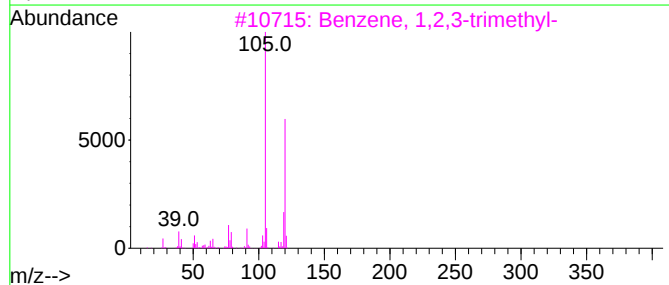
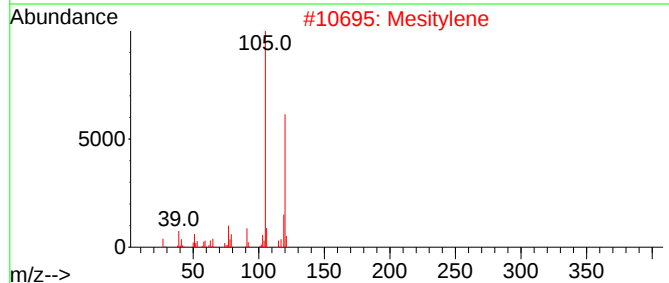
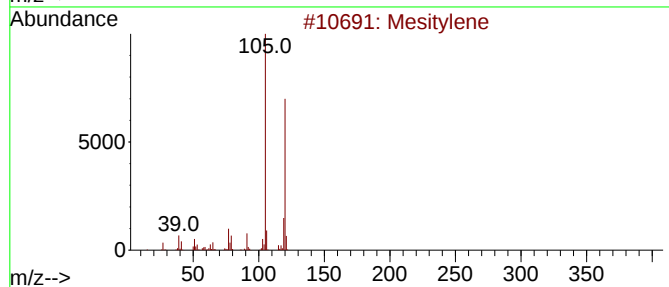
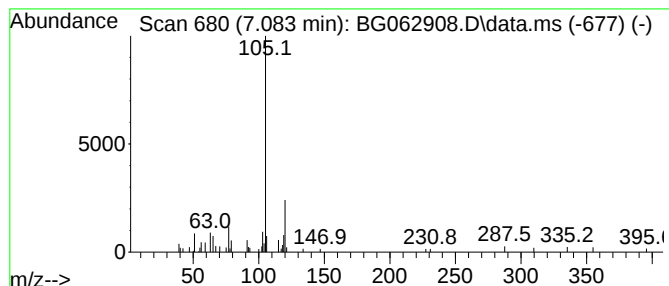
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Mesitylene Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.083	2.47 ng/ul	45833	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	74
2			Mesitylene	120	C9H12	000108-67-8	74
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	74
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	74
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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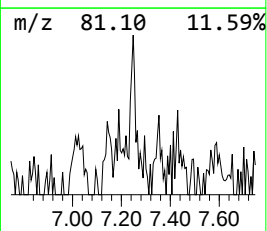
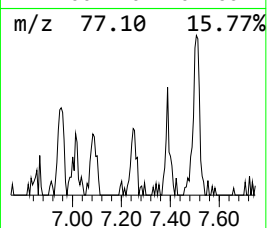
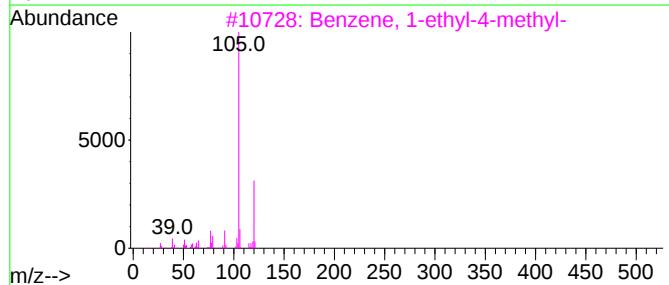
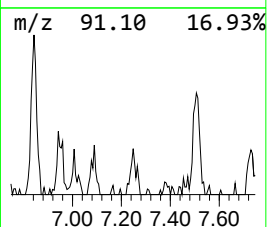
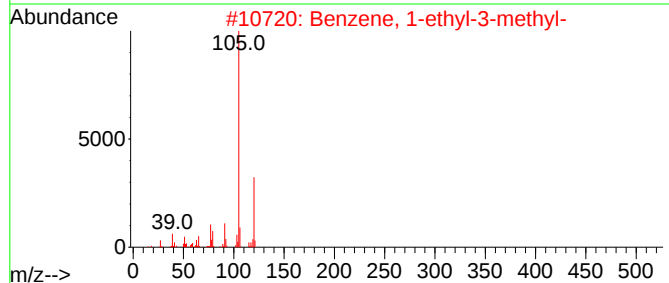
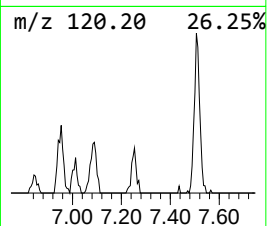
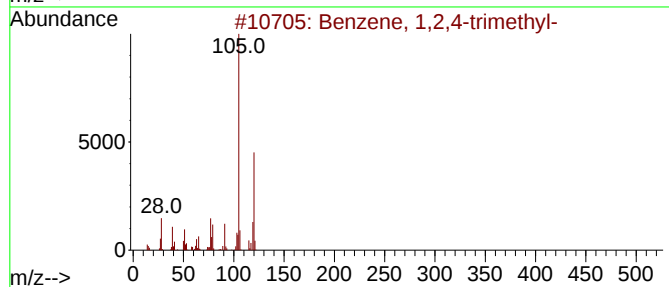
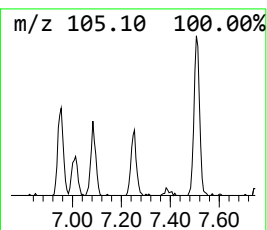
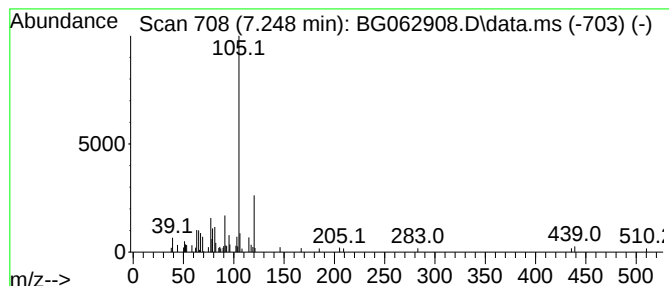
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1,2,4-trimethyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.248	2.28 ng/ul	42166	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	76
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	76
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
5			Mesitylene	120	C9H12	000108-67-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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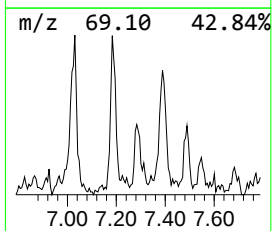
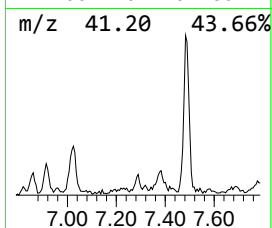
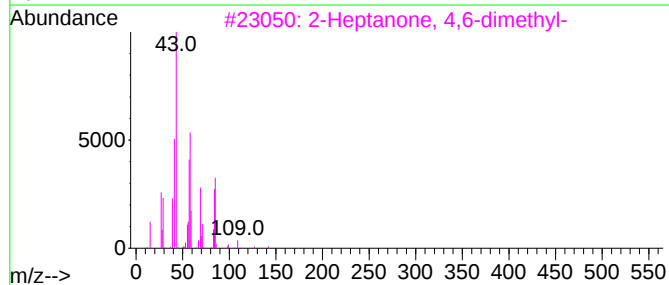
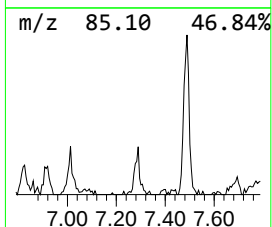
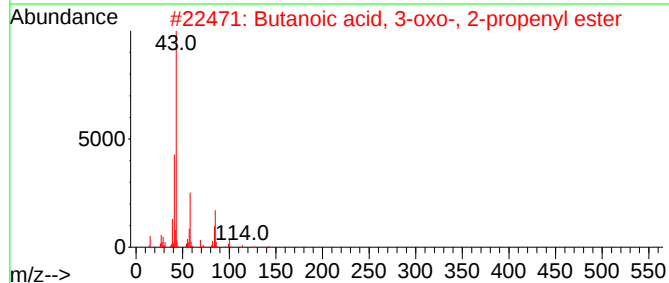
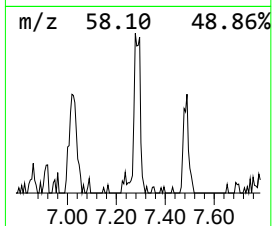
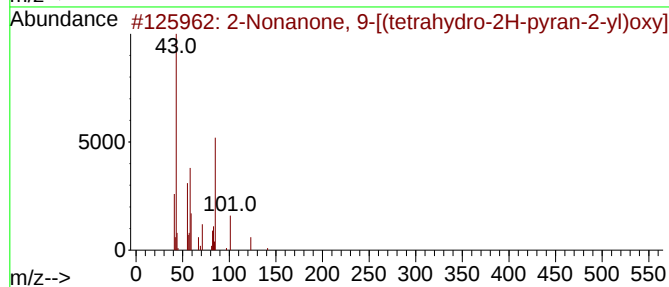
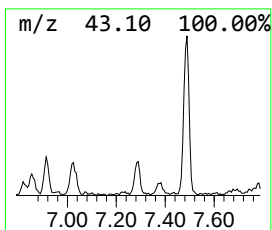
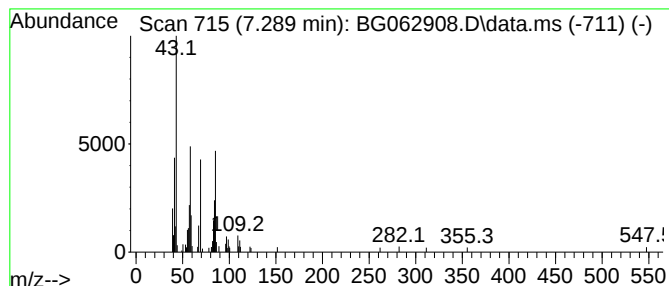
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2-Nonanone, 9-[(tetrahydro-... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.289	2.41 ng/ul	44645	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	59		
2	Butanoic acid, 3-oxo-, 2-propeny...	142	C7H10O3	001118-84-9	50		
3	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	42		
4	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	38		
5	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

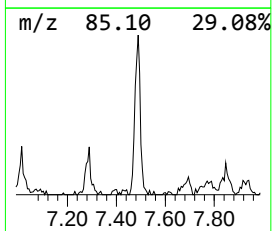
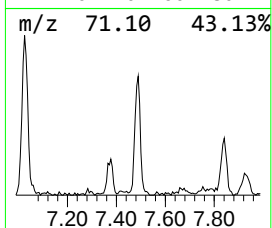
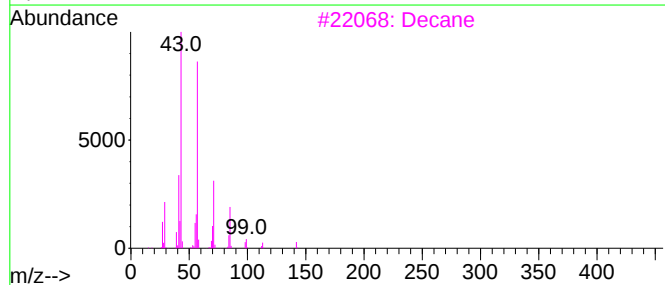
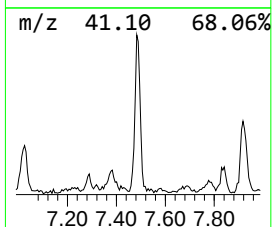
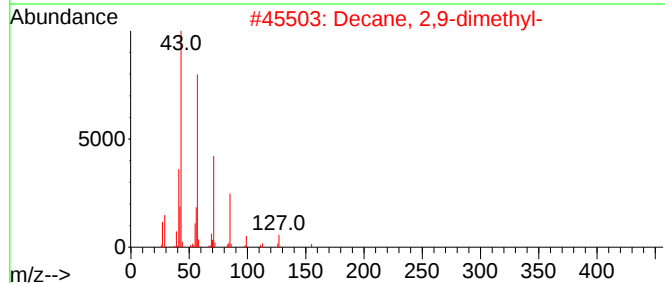
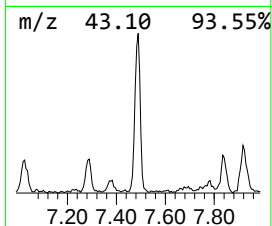
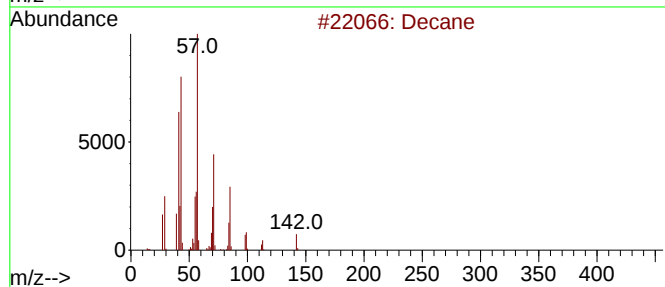
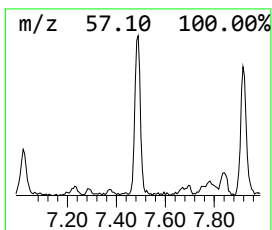
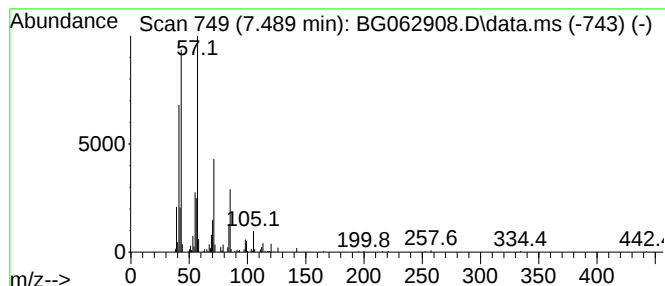
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
7.489	26.71 ng/ul	495105	1,4-Dichlorobenzene-d4		7.859	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	80
2	Decane, 2,9-dimethyl-		170	C12H26	001002-17-1	64
3	Decane		142	C10H22	000124-18-5	62
4	Undecane		156	C11H24	001120-21-4	59
5	Tridecane		184	C13H28	000629-50-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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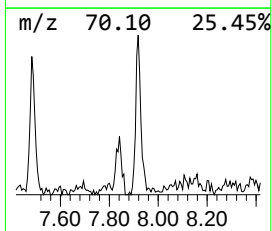
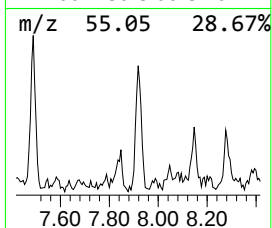
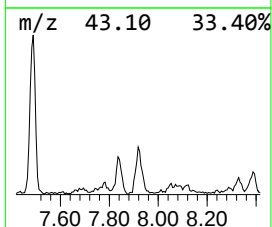
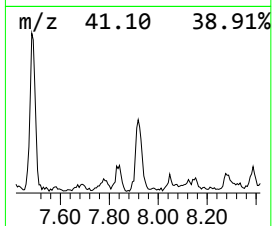
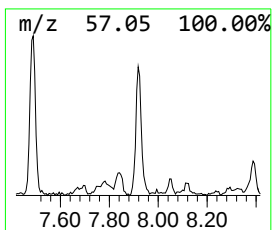
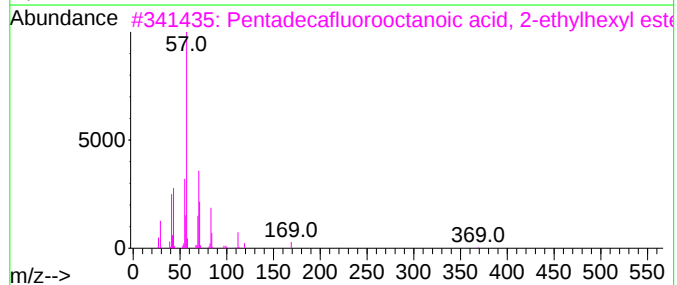
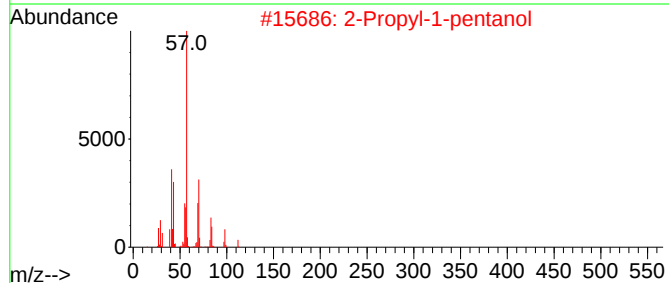
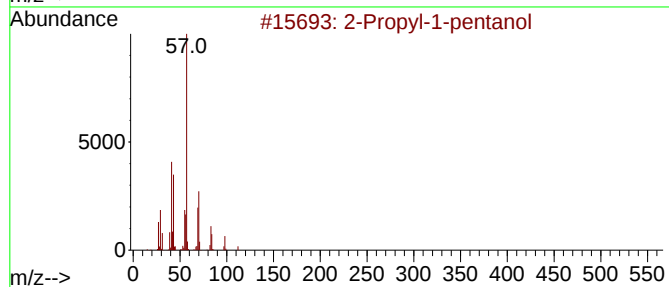
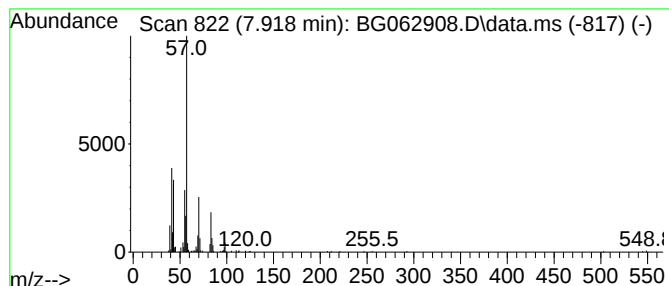
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 2-Propyl-1-pentanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.918	11.04 ng/ul	204565	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	78
2			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	56
3			Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	56
4			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	53
5			2-Ethyl-1-hexanol, trifluoroacetate	226	C10H17F3O2	053800-08-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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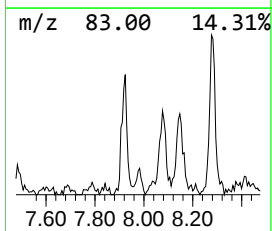
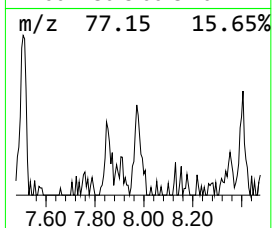
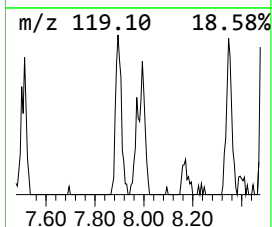
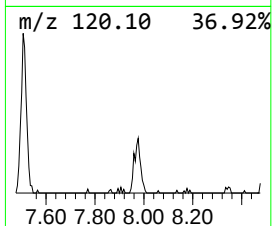
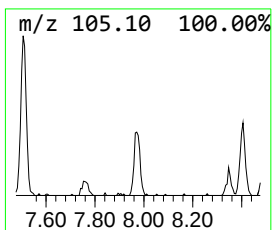
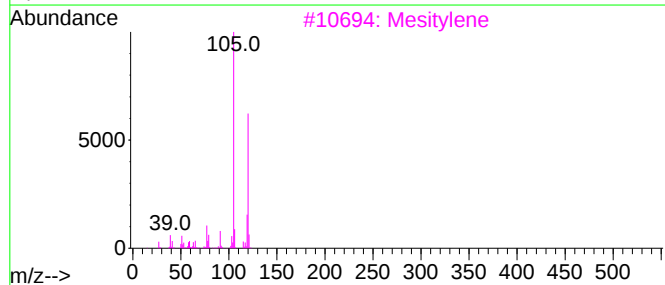
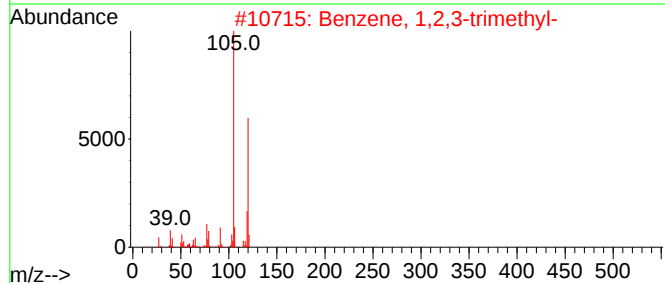
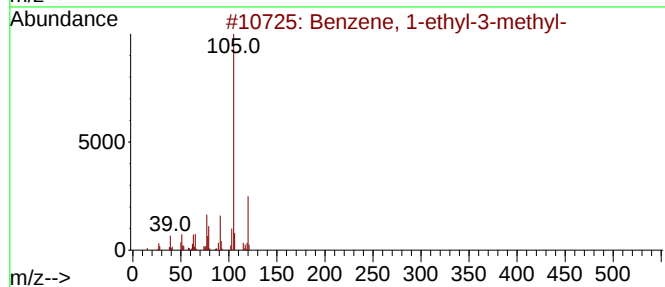
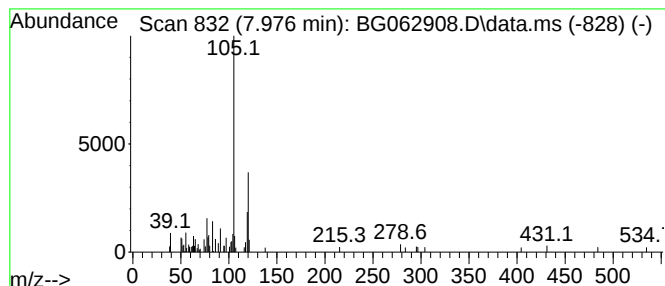
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.976	2.99 ng/ul	55450	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	64
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	62
3			Mesitylene	120	C9H12	000108-67-8	58
4			2,4-Nonadiyne	120	C9H12	063621-15-8	58
5			Mesitylene	120	C9H12	000108-67-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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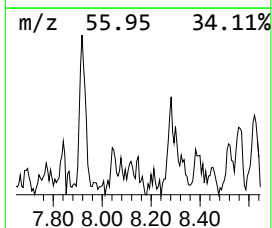
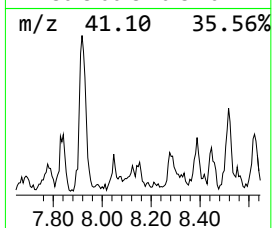
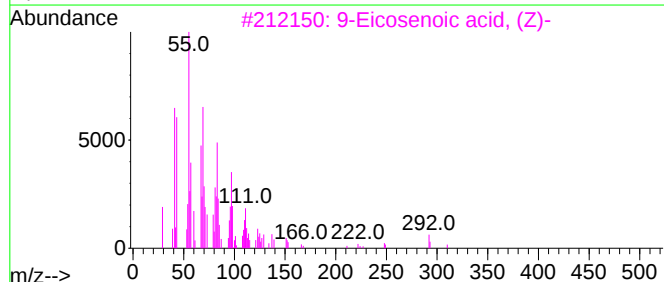
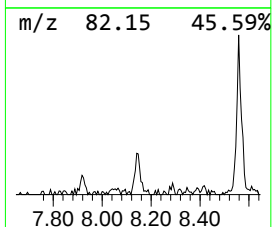
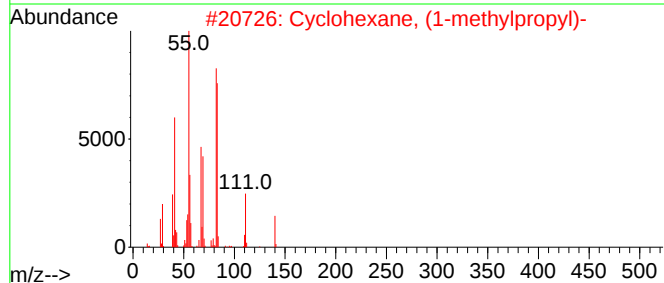
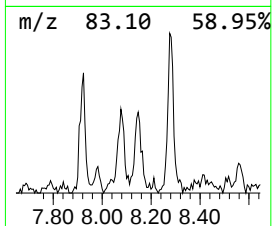
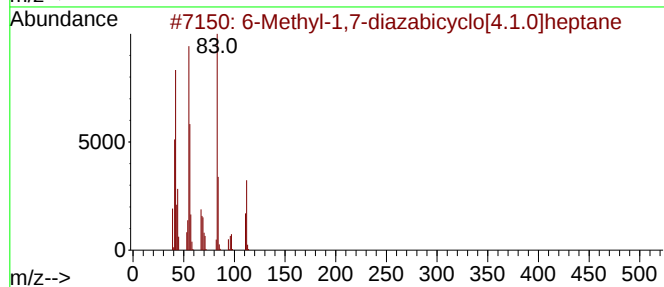
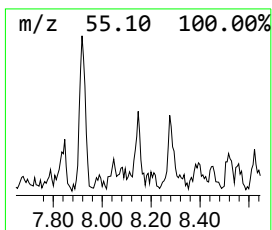
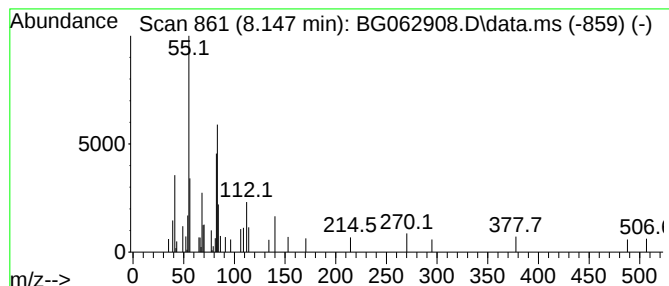
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-02 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.147	2.17 ng/ul	40242	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Methyl-1,7-diazabicyclo[4.1.0]...	112	C6H12N2	108602-71-7	43
2			Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	38
3			9-Eicosenoic acid, (Z)-	310	C20H38O2	029204-02-2	38
4			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	38
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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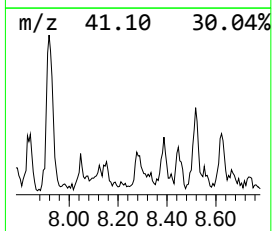
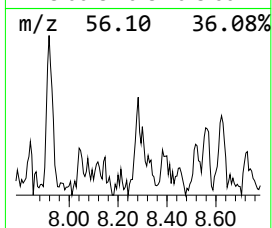
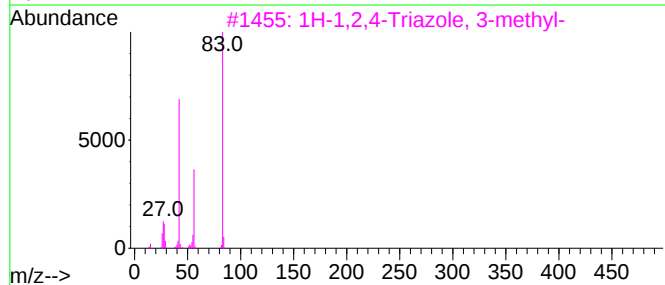
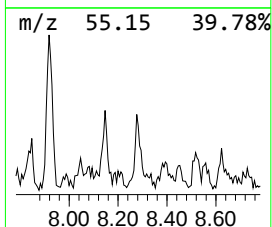
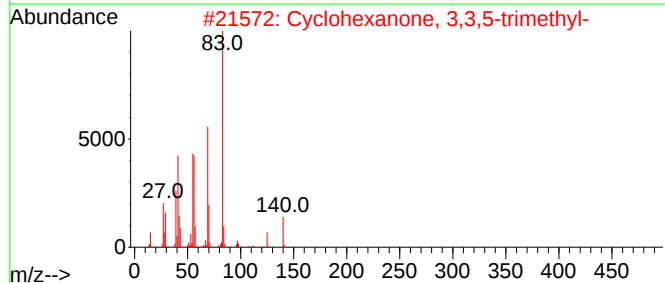
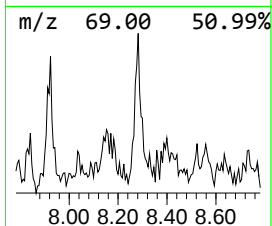
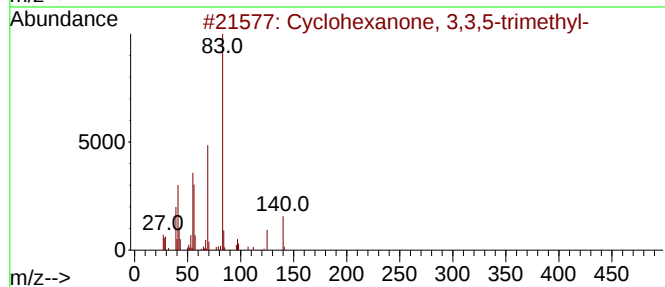
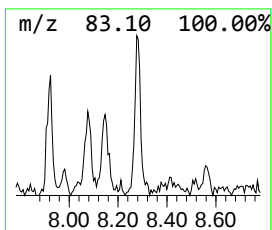
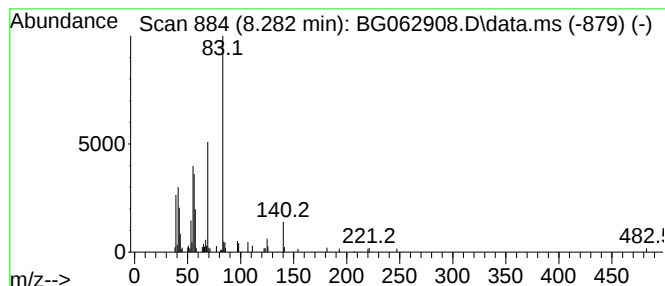
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Cyclohexanone, 3,3,5-trimet... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.282	3.76 ng/ul	69688	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	78
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	72
3			1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	64
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	59
5			Imidazole, 2-acetamido-	125	C5H7N3O	052737-49-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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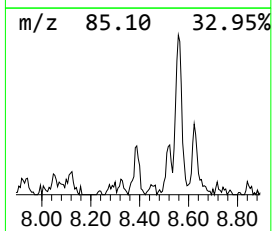
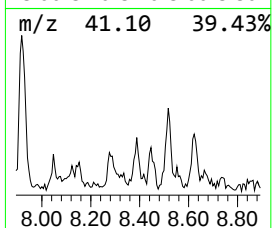
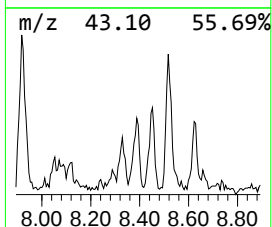
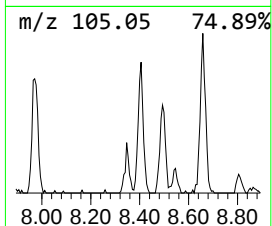
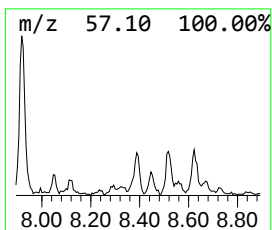
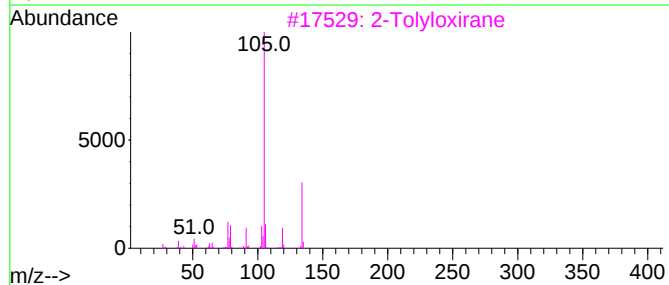
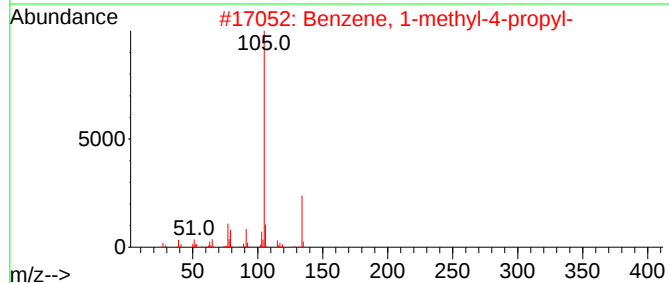
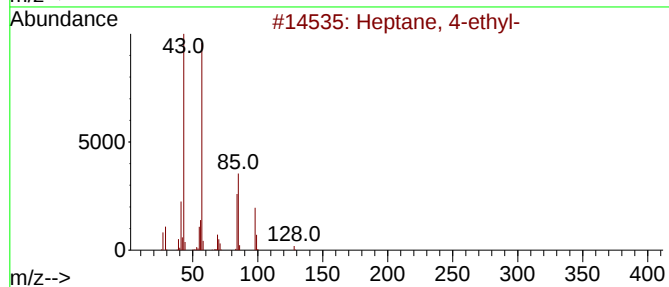
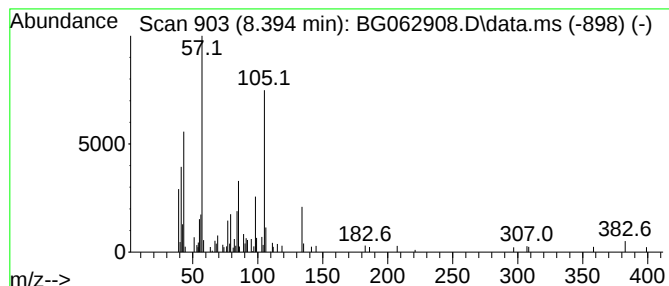
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.394	4.58 ng/ul	84810	1,4-Dichlorobenzene-d4	7.859

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-ethyl-	128	C9H20	002216-32-2	35
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	25
3		2-Tolyloxirane	134	C9H10O	002783-26-8	25
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	25
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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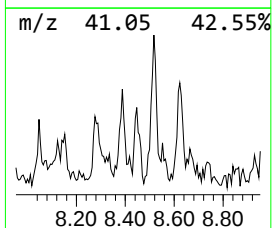
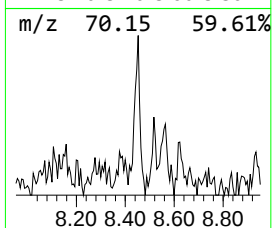
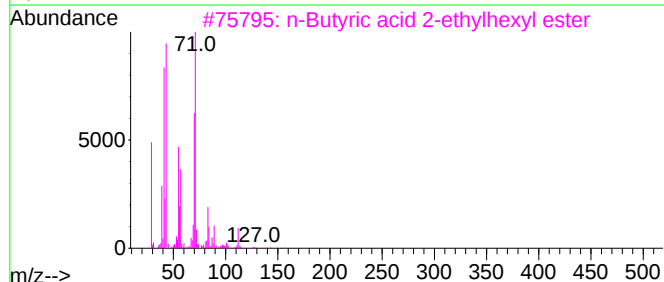
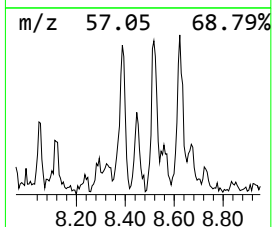
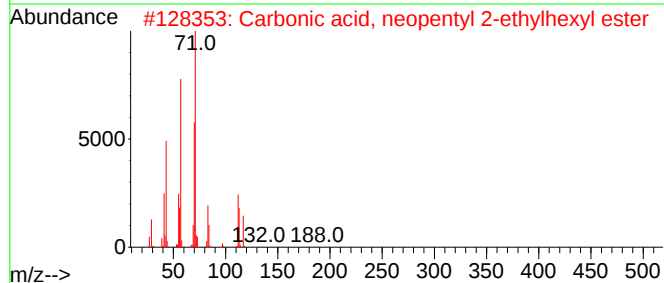
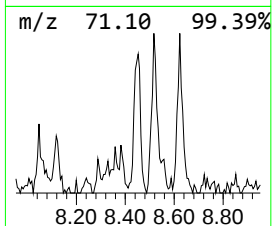
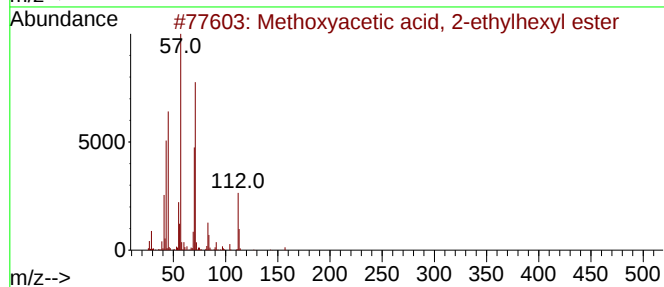
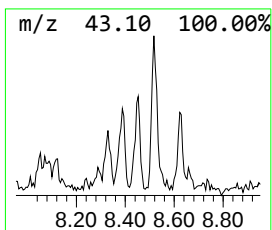
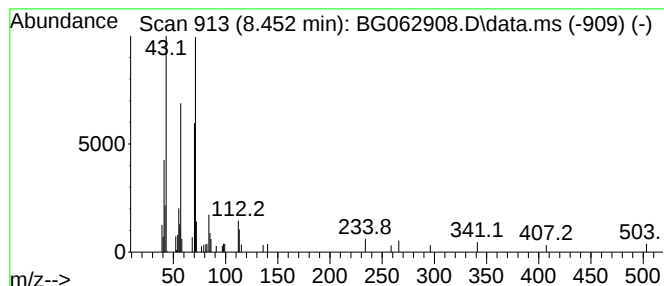
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Methoxyacetic acid, 2-ethyl... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.452	2.18 ng/ul	40340	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxyacetic acid, 2-ethylhexyl...	202	C11H22O3	1000282-41-4	72
2			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	72
3			n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	64
4			Oxalic acid, 2-ethylhexyl ethyl ...	230	C12H22O4	1000309-38-4	64
5			Decane, 4-methyl-	156	C11H24	002847-72-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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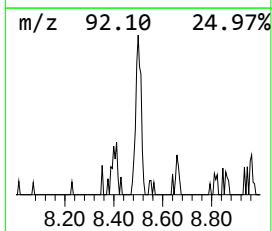
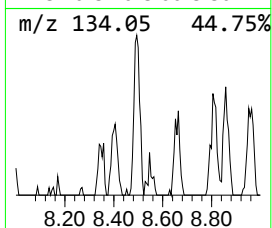
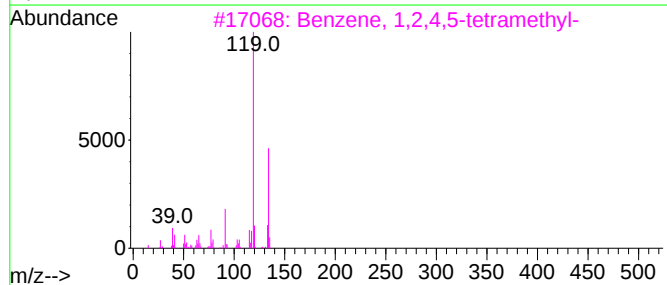
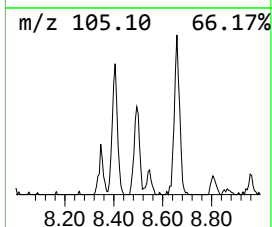
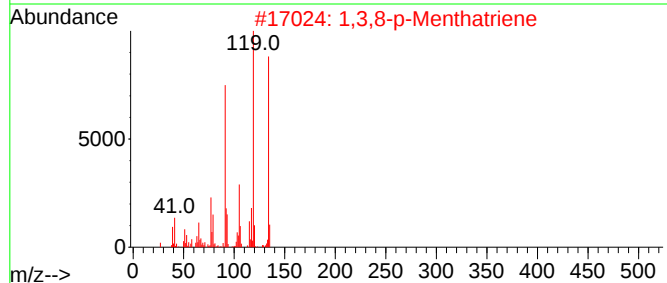
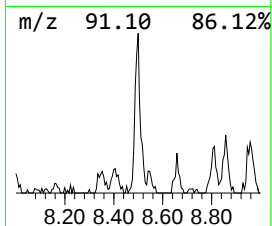
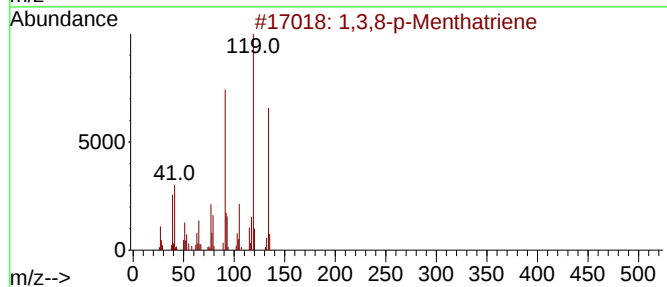
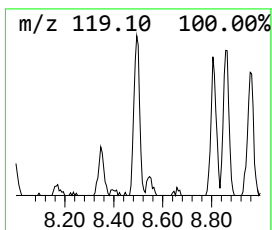
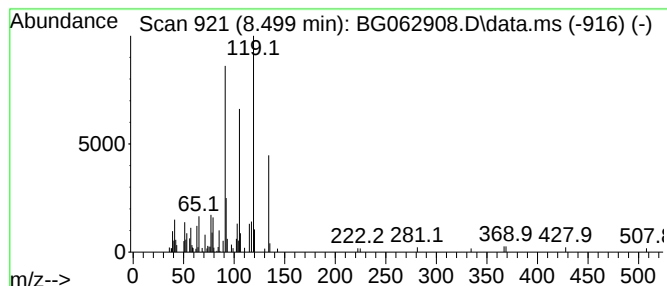
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 1,3,8-p-Menthatriene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.499	4.54 ng/ul	84161	1,4-Dichlorobenzene-d4	7.859

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	93
2		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	87
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87
4		1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	87
5		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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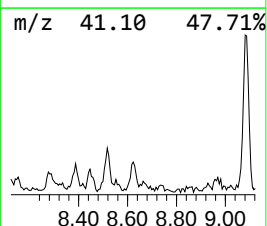
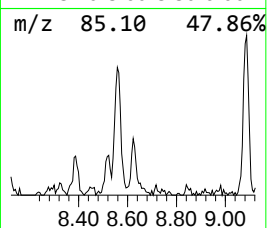
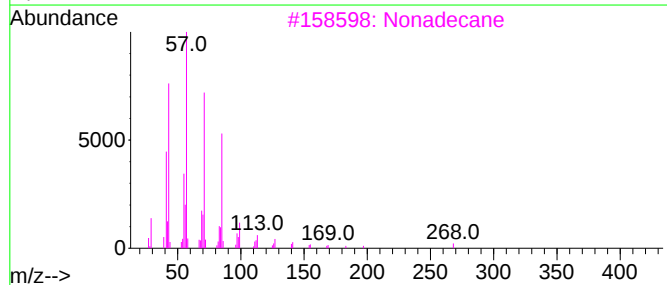
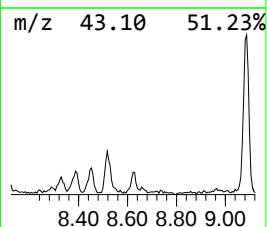
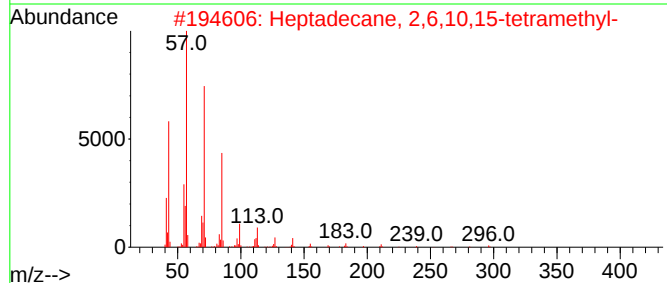
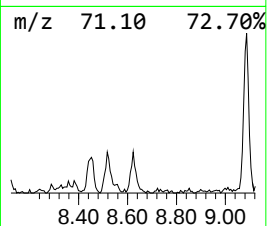
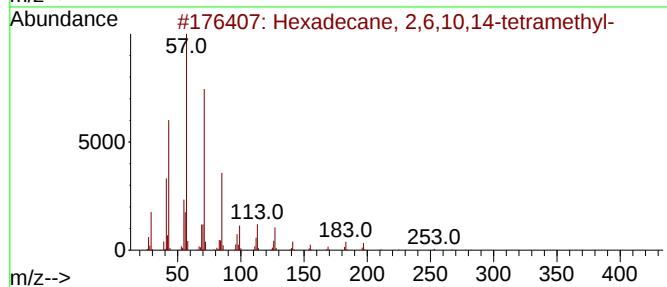
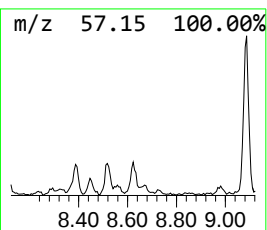
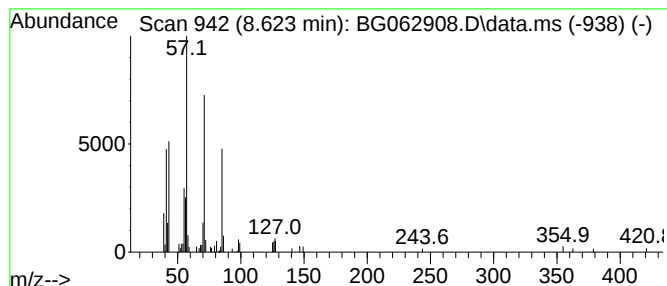
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.623	3.68 ng/ul	68129	1,4-Dichlorobenzene-d4	7.859

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
3		Nonadecane	268	C19H40	000629-92-5	72
4		Hexadecane	226	C16H34	000544-76-3	72
5		Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCVZ0ME

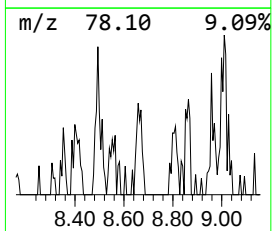
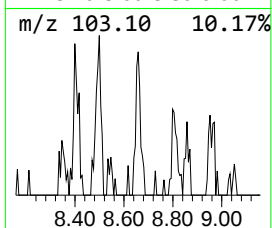
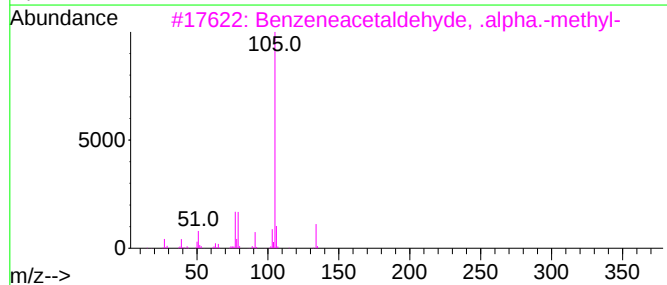
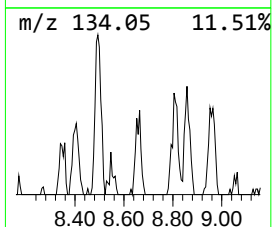
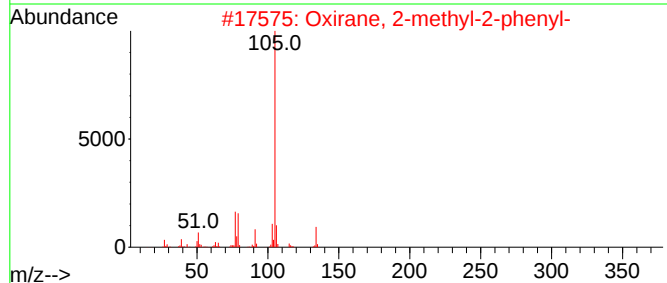
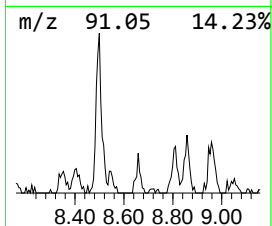
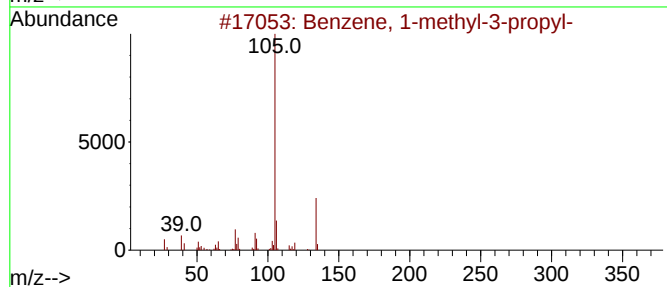
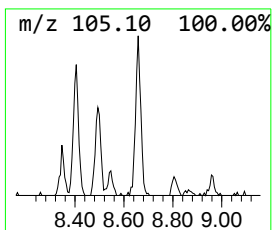
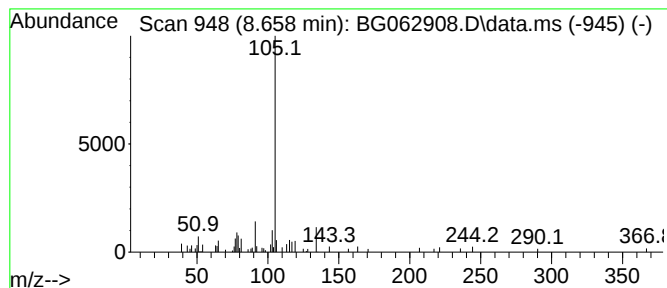
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 1-methyl-3-propyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.658	3.64 ng/ul	67490	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	50
2			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	49
3			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	47
4			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	47
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

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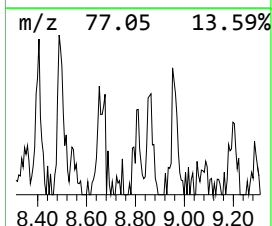
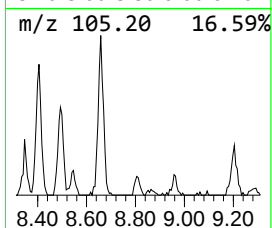
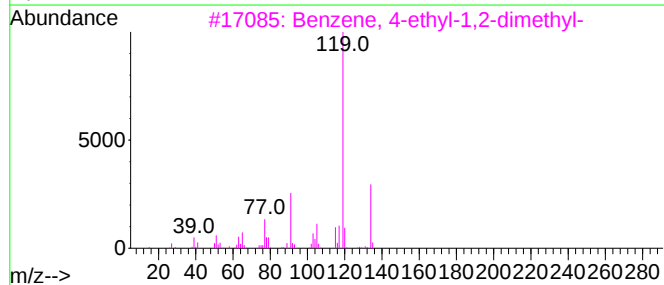
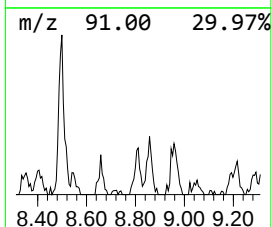
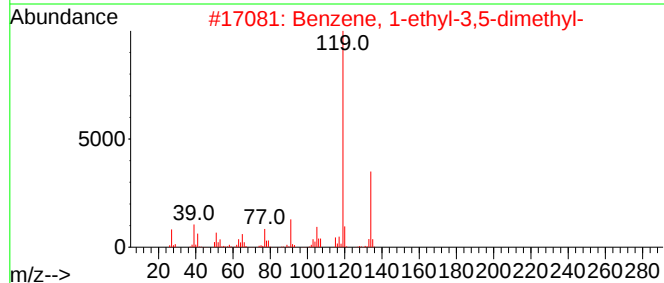
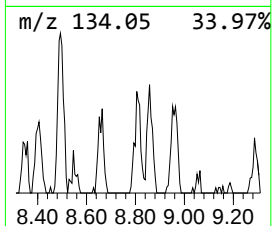
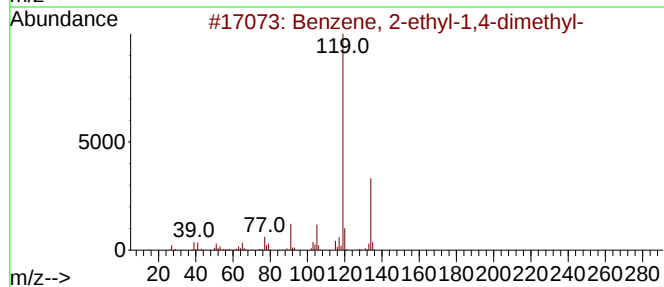
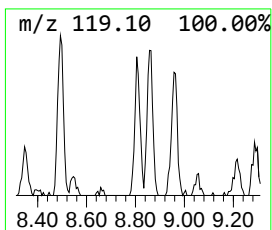
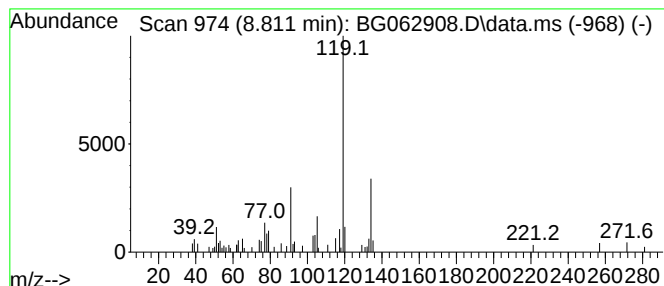
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.811	2.60 ng/ul	48231	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

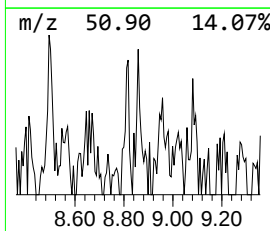
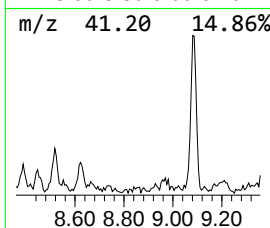
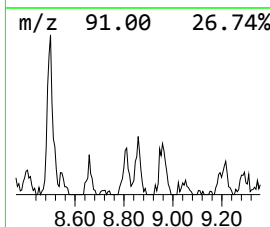
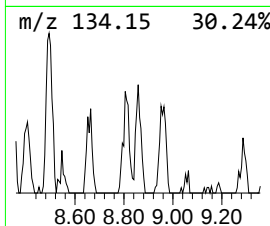
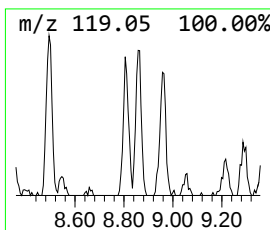
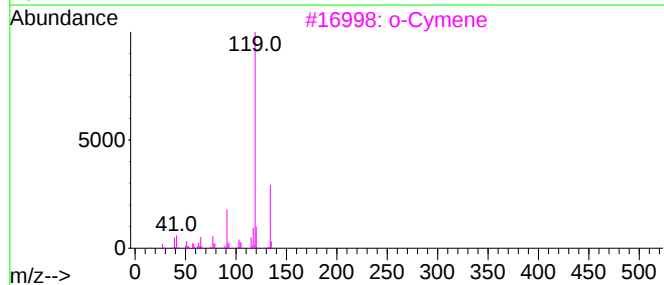
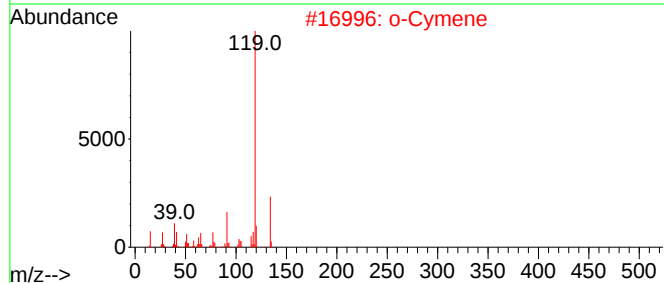
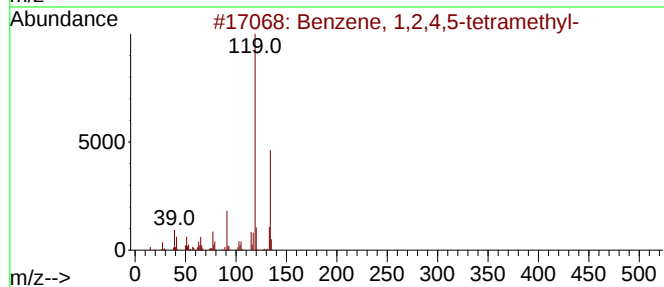
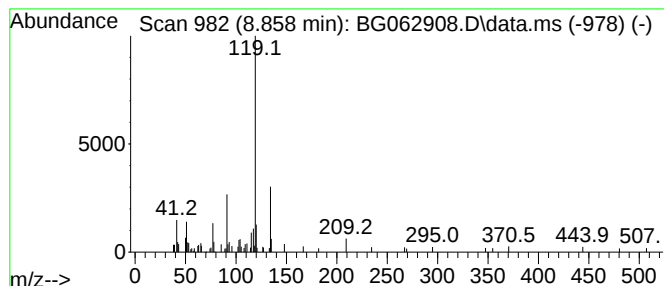
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.858	3.11 ng/ul	57601	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87
2			o-Cymene	134	C10H14	000527-84-4	87
3			o-Cymene	134	C10H14	000527-84-4	87
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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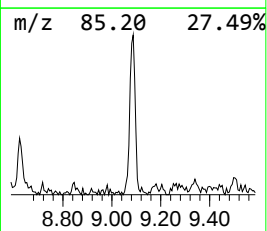
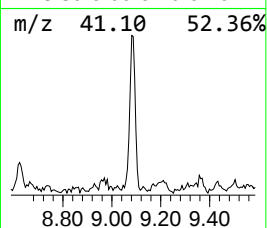
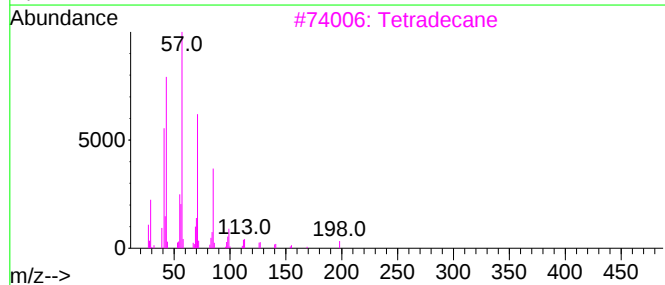
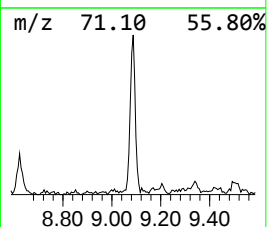
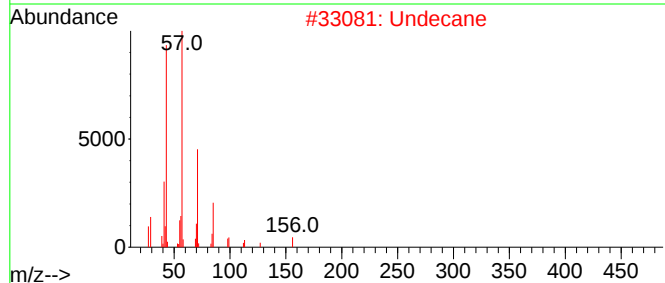
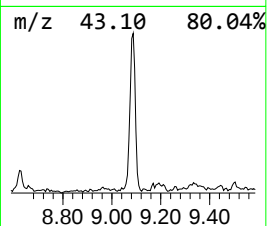
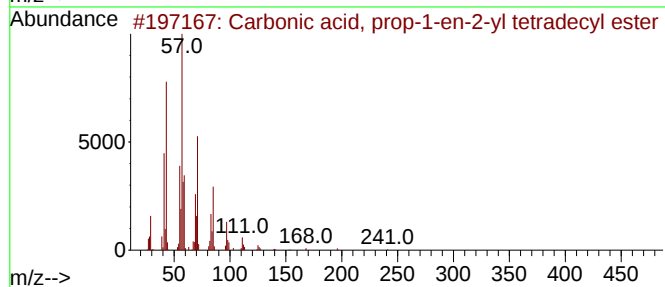
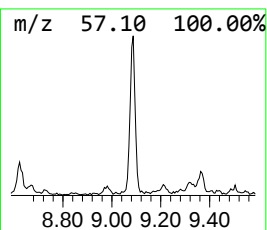
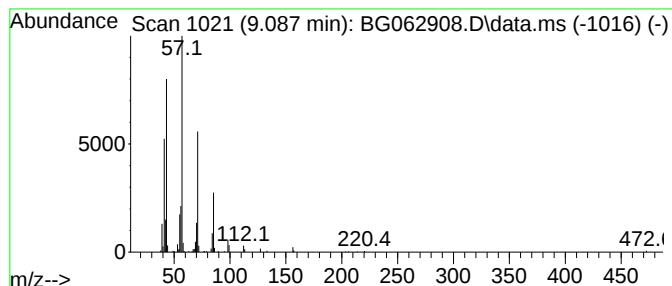
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Carbonic acid, prop-1-en-2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.087	20.00 ng/ul	370695	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
2			Undecane	156	C11H24	001120-21-4	86
3			Tetradecane	198	C14H30	000629-59-4	83
4			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
5			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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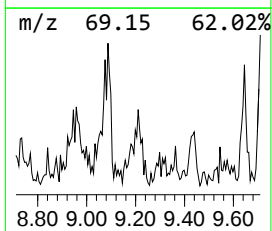
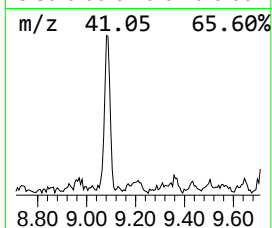
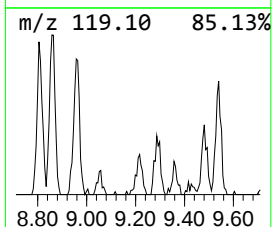
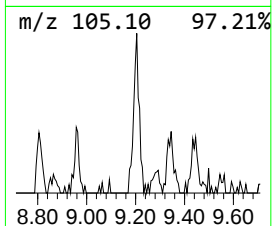
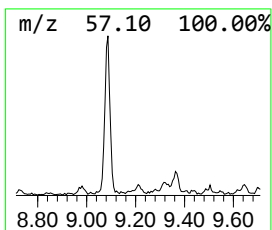
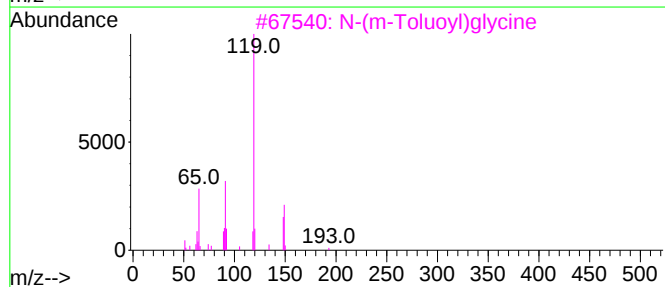
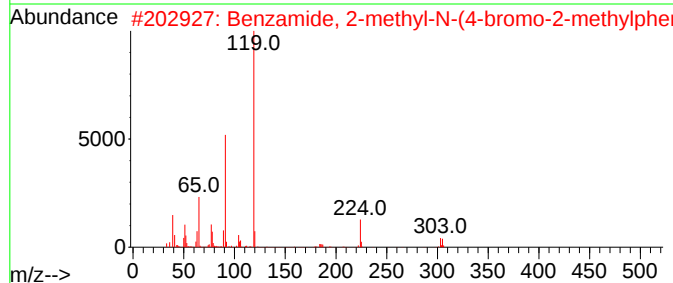
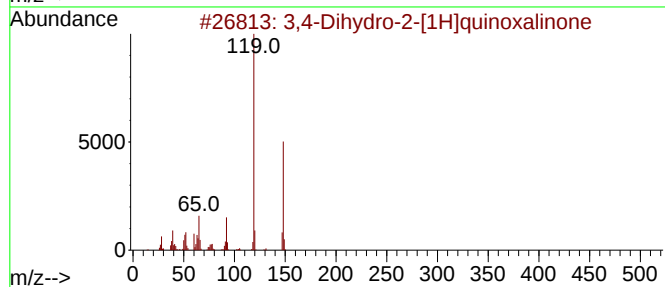
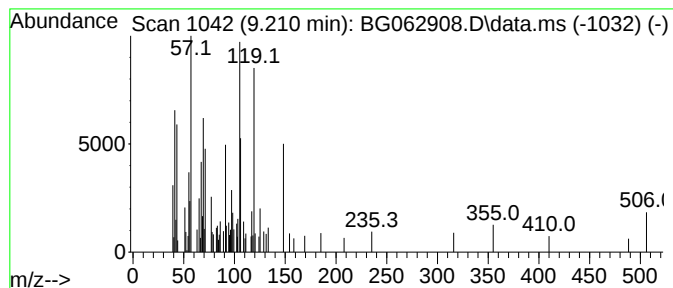
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.210	5.66 ng/ul	104863	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dihydro-2-[1H]quinoxalinone	148	C8H8N2O	059564-59-9	38
2			Benzamide, 2-methyl-N-(4-bromo-2...	303	C15H14BrNO	306743-63-5	22
3			N-(m-Toluoyl)glycine	193	C10H11NO3	027115-49-7	22
4			5-Aza-2H-indazole	119	C6H5N3	000271-50-1	18
5			3-Toluoylacetonitrile	159	C10H9NO	053882-81-8	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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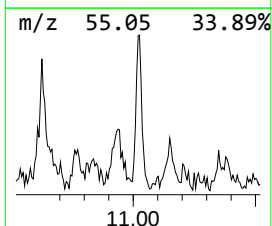
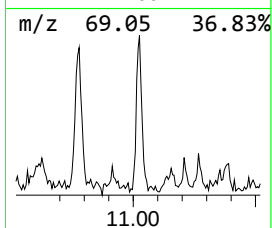
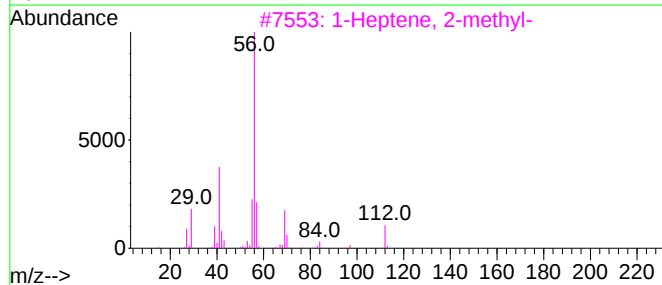
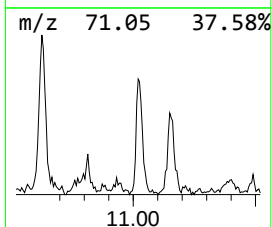
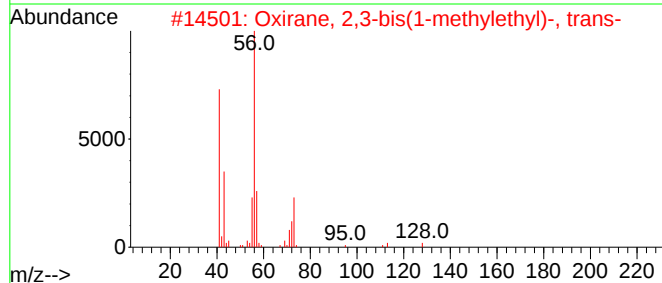
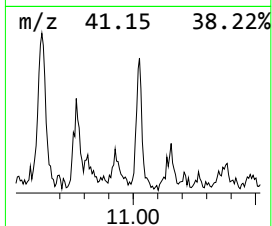
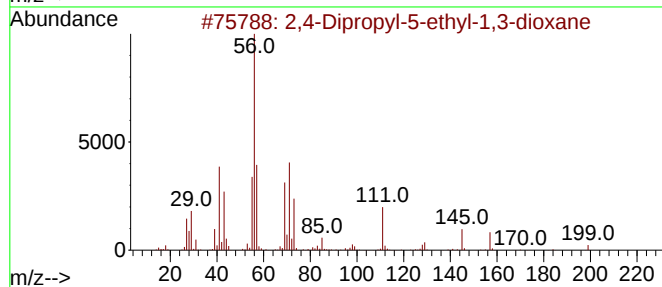
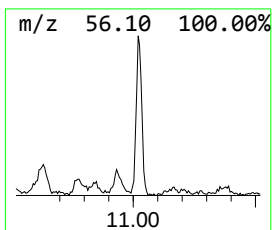
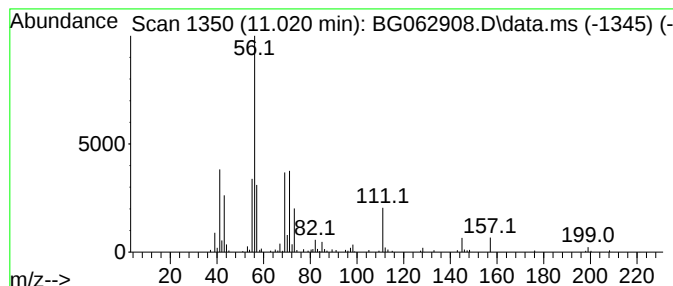
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.020	5.75 ng/ul	221975	Naphthalene-d8	10.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	86	
2	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	38	
3	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	38	
4	2-Methyl-1-nonene	140	C10H20	002980-71-4	35	
5	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	33	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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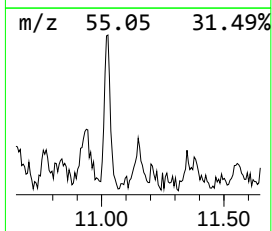
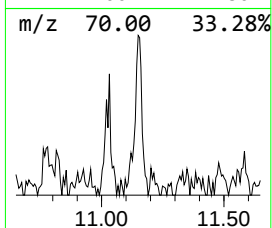
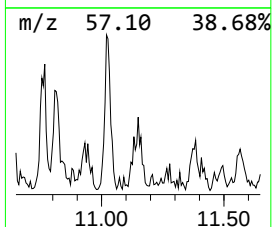
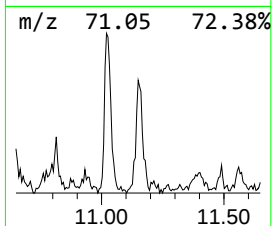
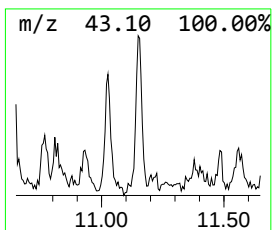
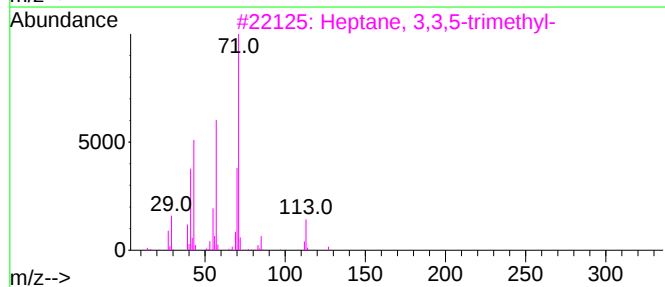
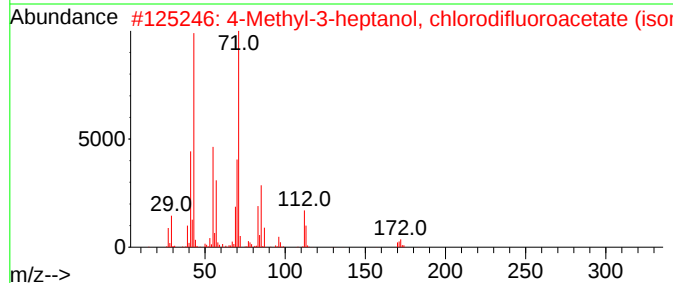
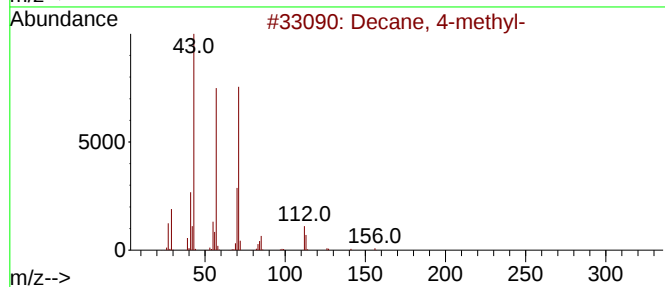
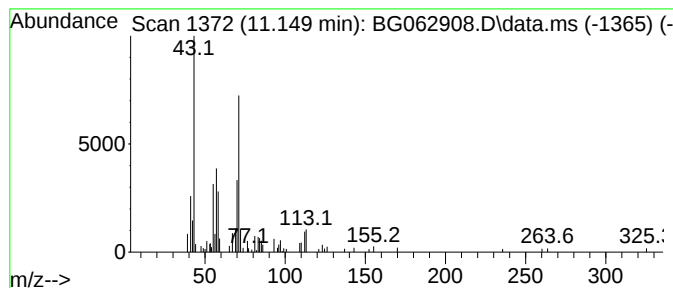
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.149	2.89 ng/ul	111453	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	53
2			4-Methyl-3-heptanol, chlorodiflu...	242	C10H17ClF2O2	1000447-28-3	50
3			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	50
4			Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	50
5			Tetradecane, 4-methyl-	212	C15H32	025117-24-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
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Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

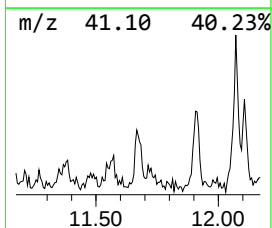
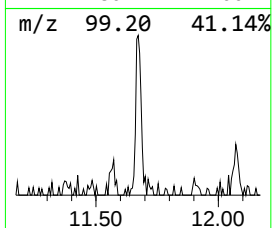
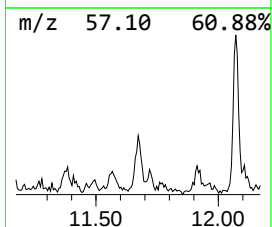
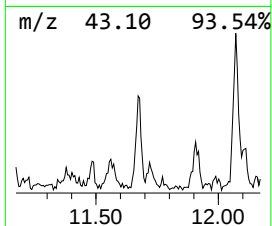
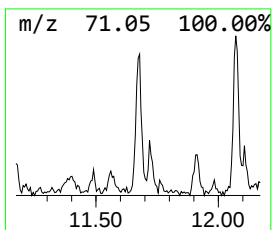
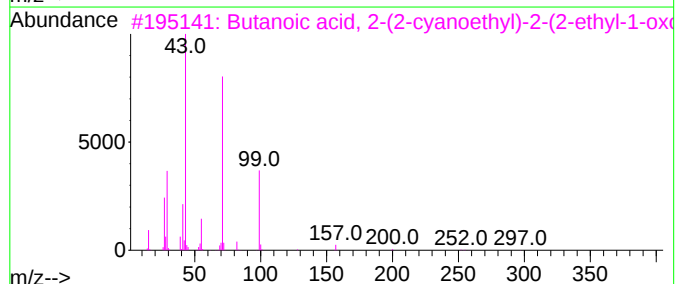
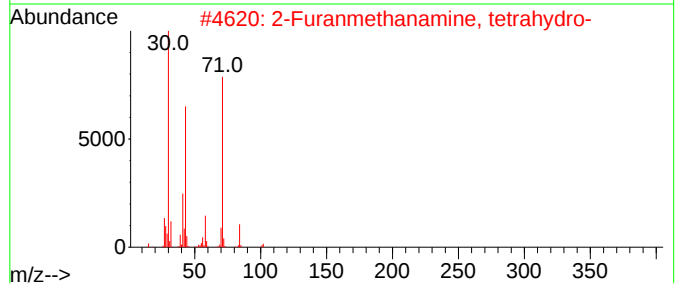
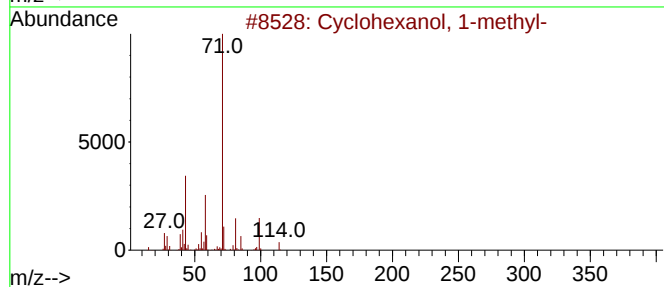
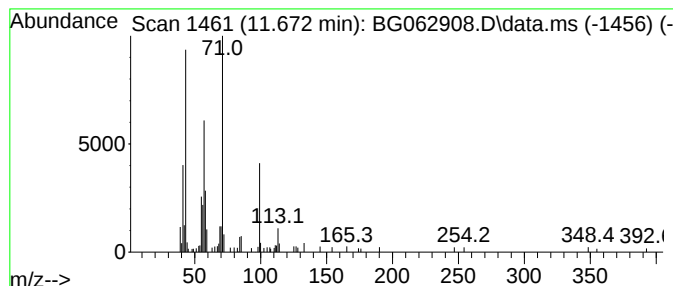
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Cyclohexanol, 1-methyl- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.672	2.58 ng/ul	99522	Naphthalene-d8			10.644
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexanol, 1-methyl-		114	C7H14O	000590-67-0	52
2	2-Furanmethanamine, tetrahydro-		101	C5H11NO	004795-29-3	47
3	Butanoic acid, 2-(2-cyanoethyl)-...		297	C15H23NO5	1010460-88-7	47
4	2-Ethylbutyric acid, pentafluoro...		282	C12H11F5O2	1010370-22-8	47
5	4-Nonanone		142	C9H18O	004485-09-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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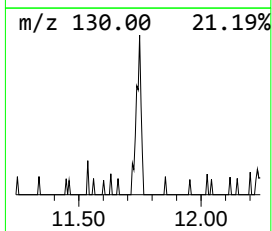
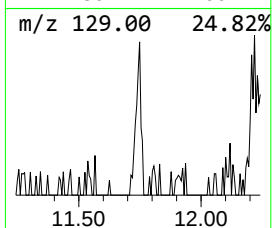
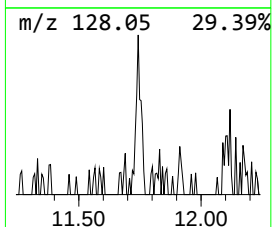
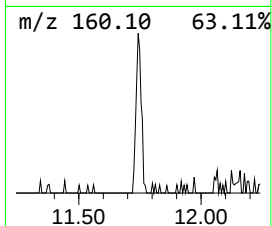
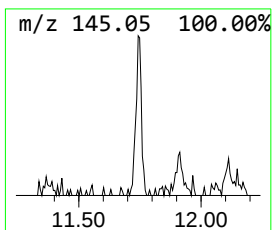
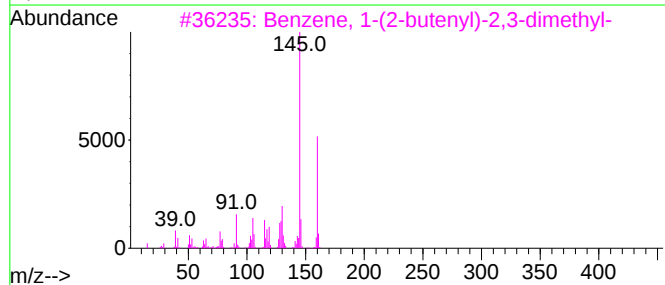
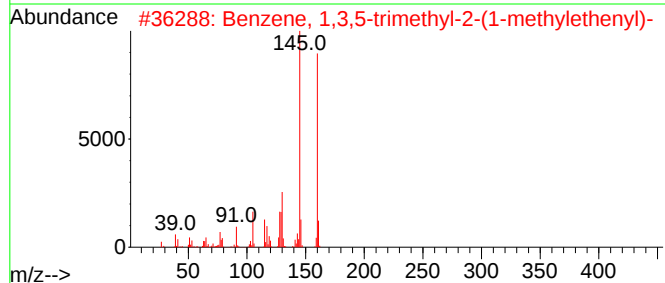
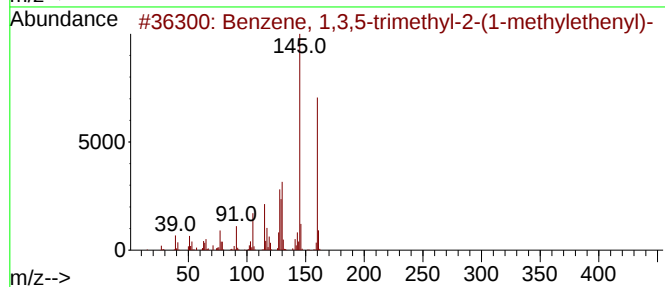
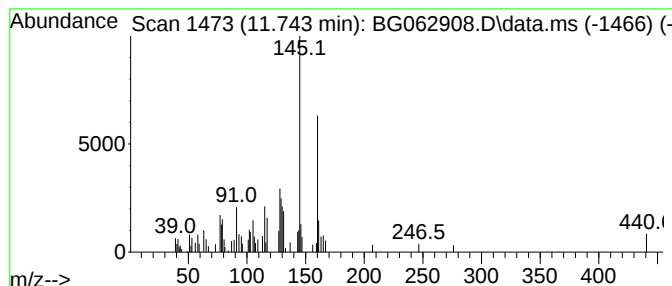
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.743	2.25 ng/ul	86842	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	87
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	87
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	74
4			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	72
5			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

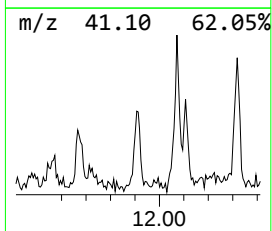
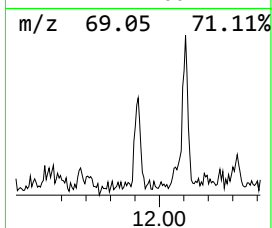
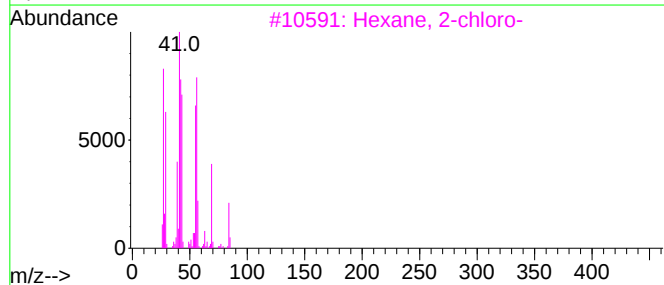
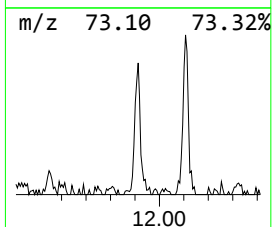
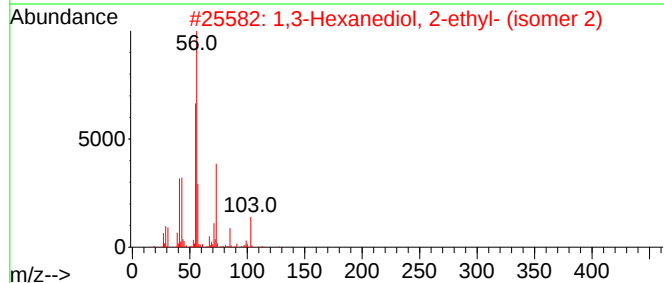
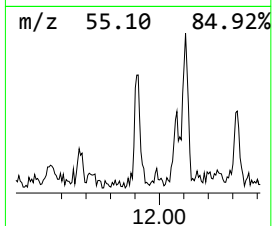
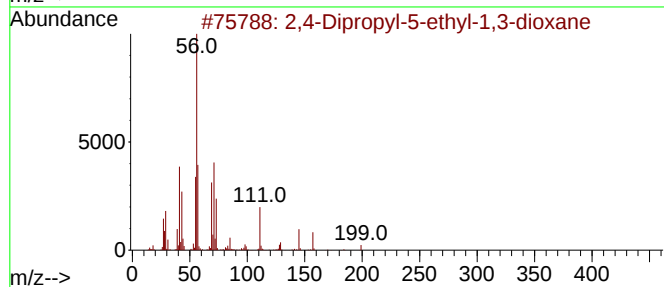
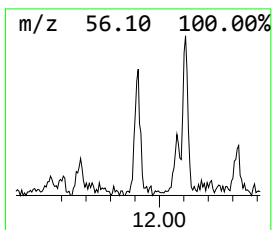
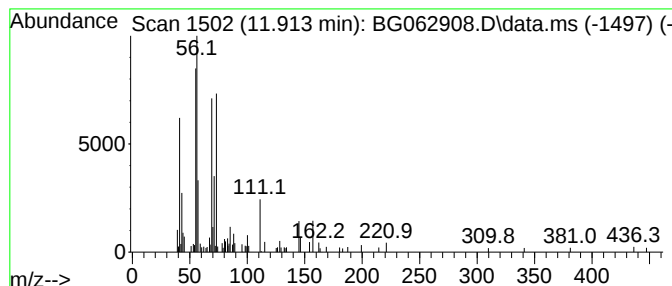
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 1,3-Hexanediol, 2-ethyl- (i... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.913	2.96 ng/ul	114182	Naphthalene-d8	10.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	56	
2	1,3-Hexanediol, 2-ethyl- (isomer 2)	146	C8H18O2	1000484-18-1	53	
3	Hexane, 2-chloro-	120	C6H13Cl	000638-28-8	43	
4	1-Decene, 2-methyl-	154	C11H22	013151-27-4	38	
5	2-Decene, 7-methyl-, (Z)-	154	C11H22	074630-23-2	38	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

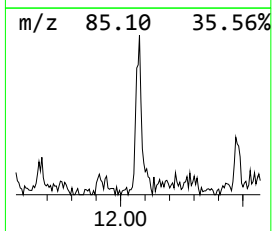
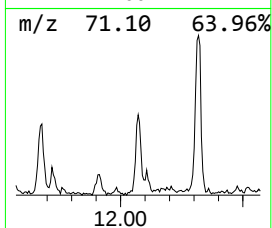
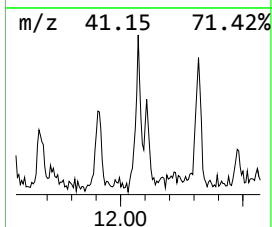
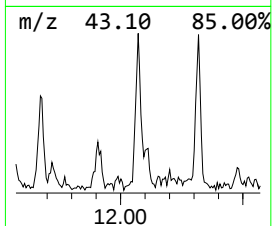
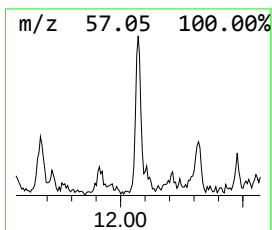
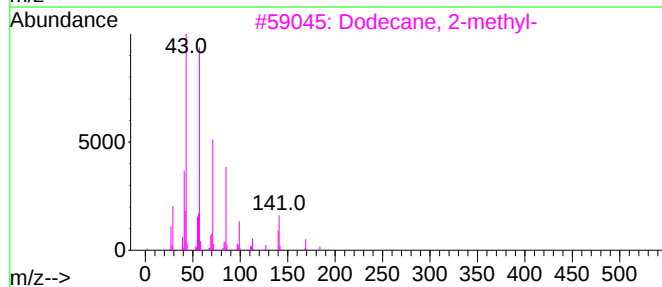
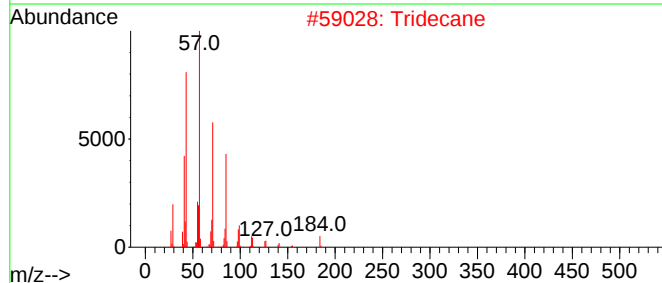
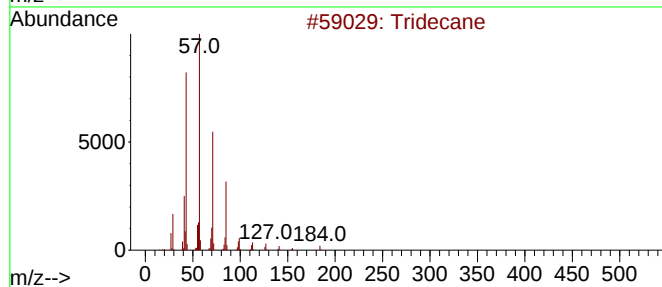
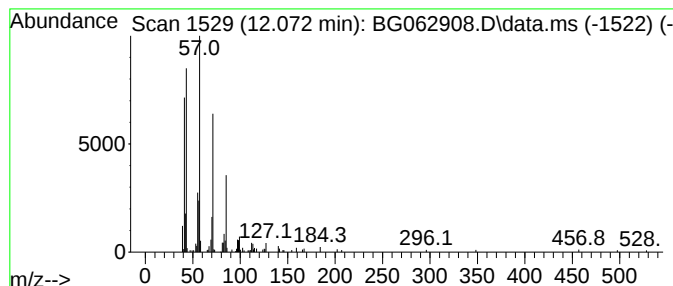
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.072	4.53 ng/ul	174875	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	94
2	Tridecane	184	C13H28	000629-50-5	74
3	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

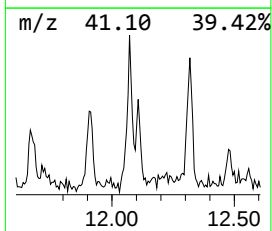
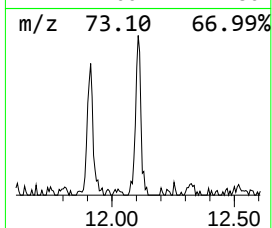
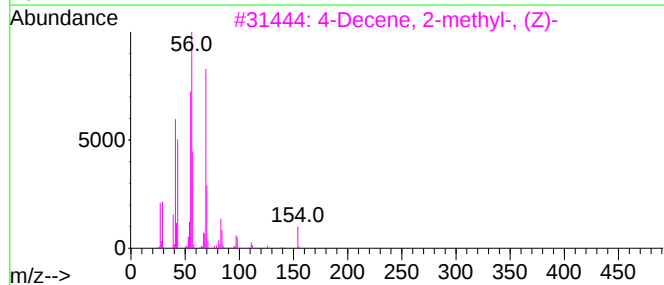
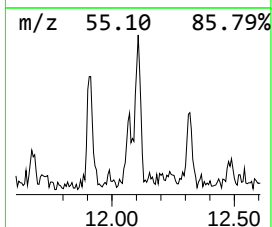
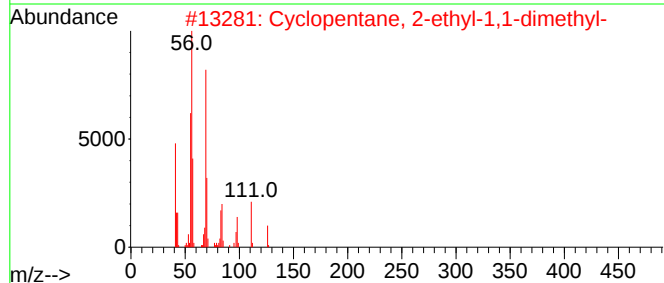
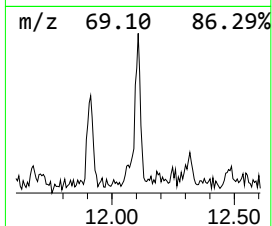
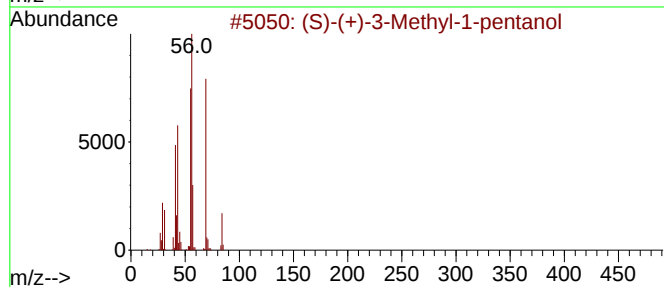
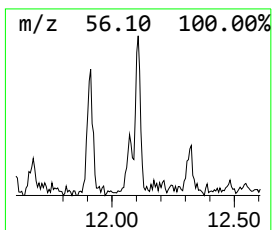
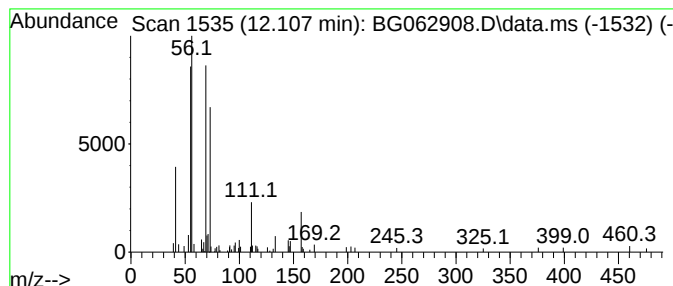
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 unknown-04 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.107	2.88 ng/ul	111294	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-3-Methyl-1-pentanol	102	C6H14O	042072-39-9	45
2	Cyclopentane, 2-ethyl-1,1-dimethyl-	126	C9H18	054549-80-3	40
3	4-Decene, 2-methyl-, (Z)-	154	C11H22	055499-07-5	38
4	Undecanol-4	172	C11H24O	004272-06-4	38
5	1-Octene, 2,7-dimethyl-	140	C10H20	033718-03-5	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

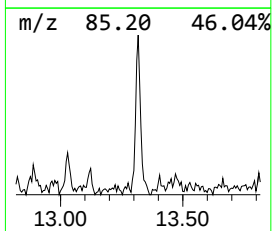
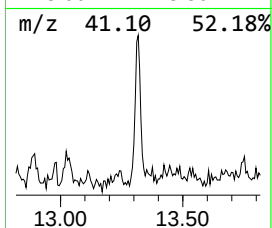
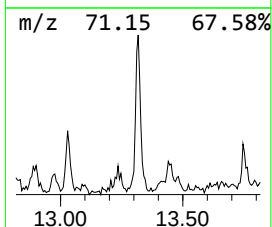
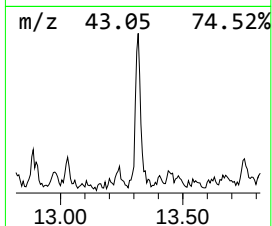
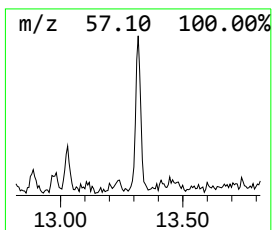
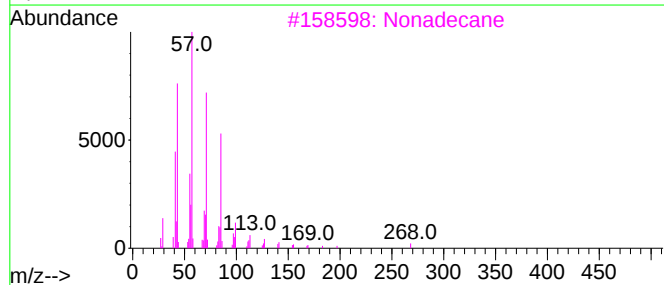
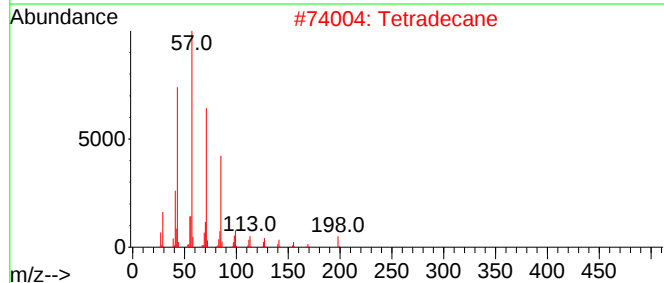
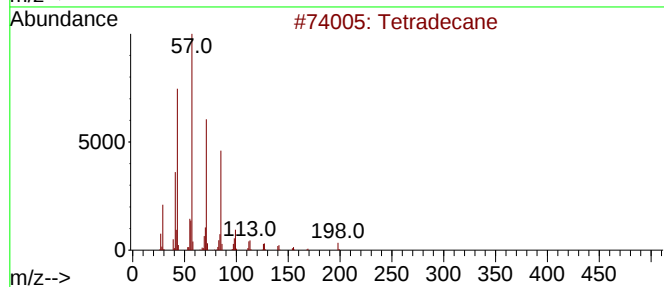
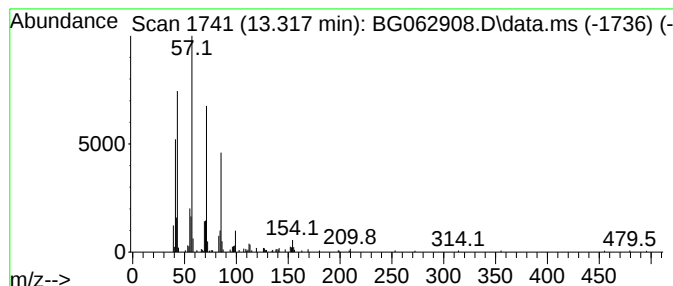
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.317	5.41 ng/ul	204750	Acenaphthene-d10		14.486	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	90
2	Tetradecane		198	C14H30	000629-59-4	86
3	Nonadecane		268	C19H40	000629-92-5	86
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	78
5	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

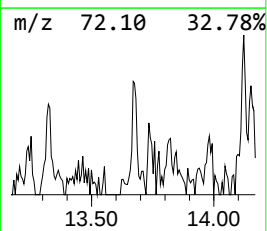
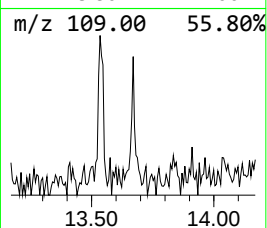
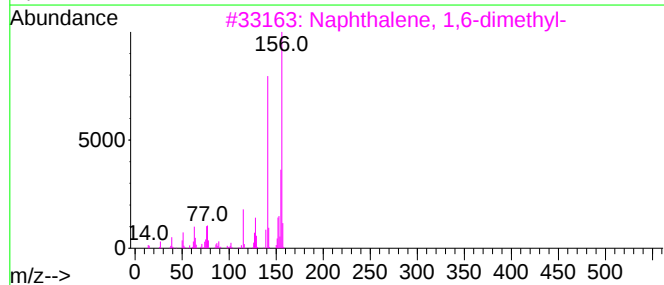
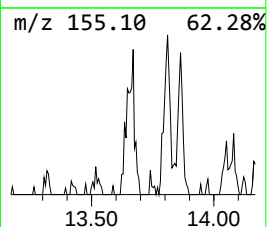
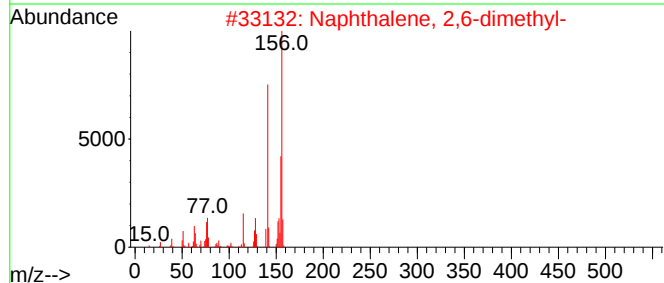
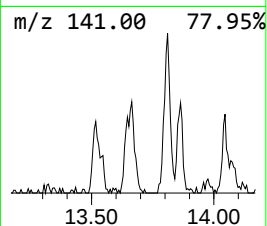
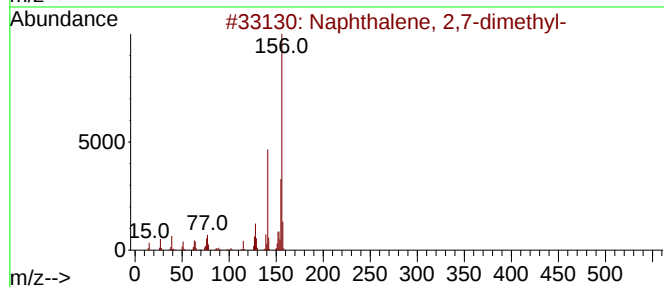
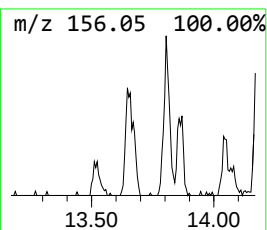
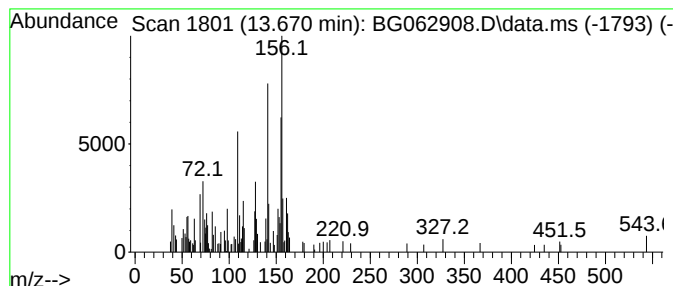
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Naphthalene, 2,7-dimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.670	2.74 ng/ul	103606	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	60
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	52
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	49
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	49
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

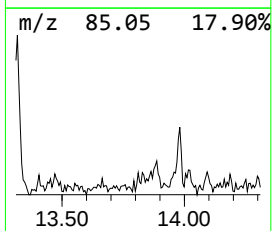
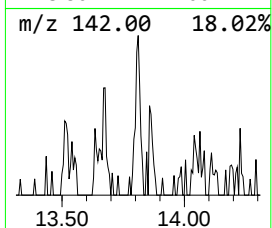
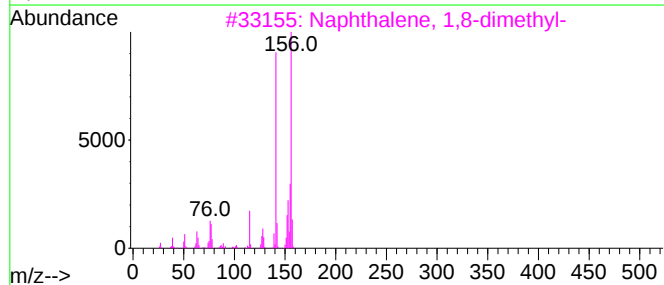
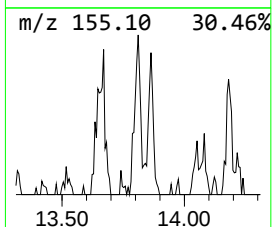
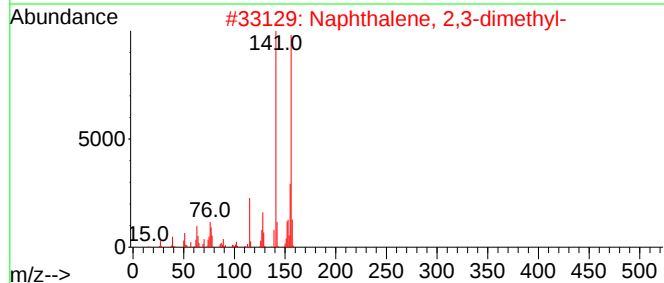
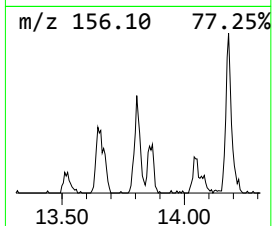
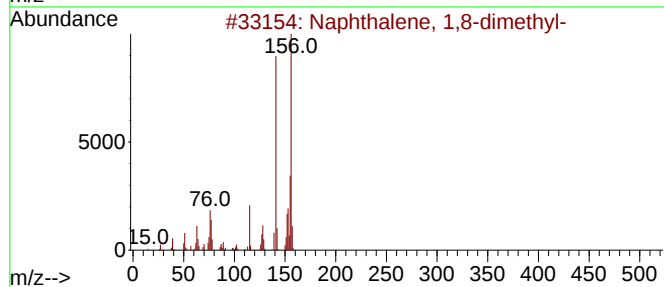
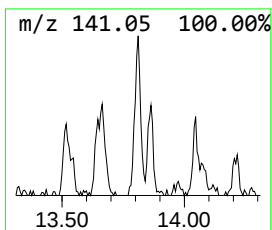
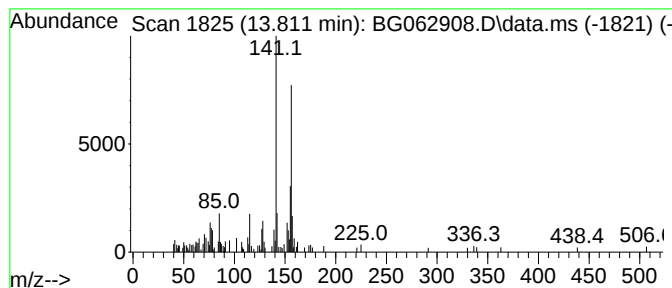
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Naphthalene, 1,8-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.811	3.05 ng/ul	115283	Acenaphthene-d10	14.486

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
5		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

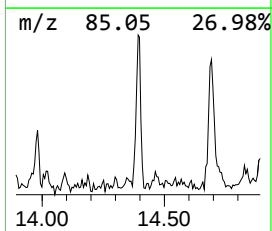
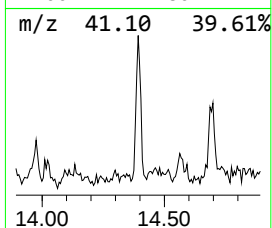
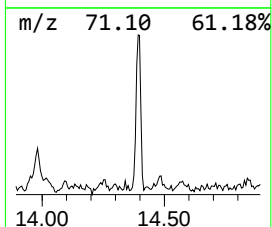
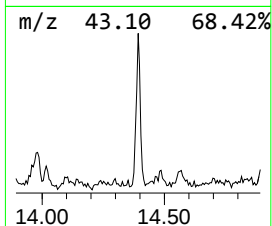
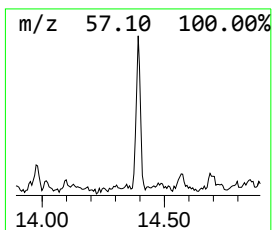
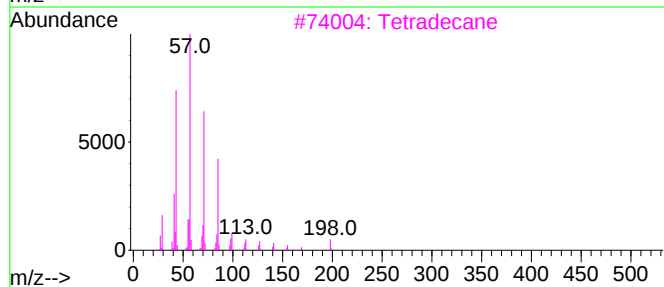
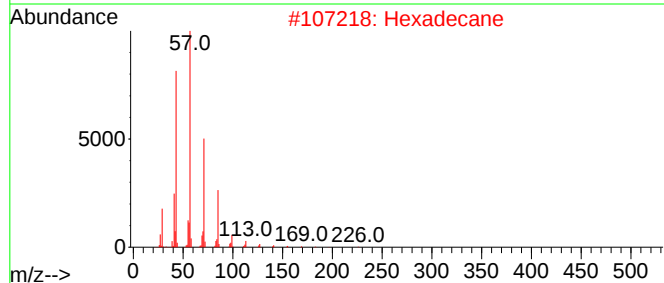
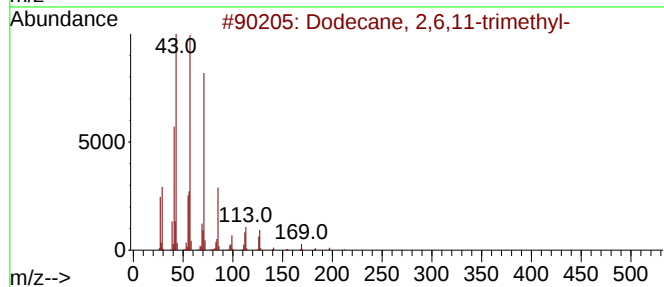
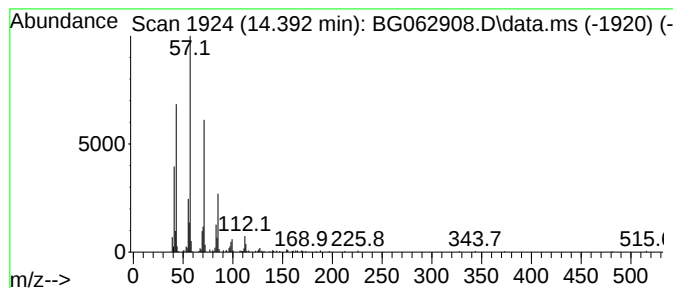
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.392	6.31 ng/ul	238515	Acenaphthene-d10	14.486

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	91
2		Hexadecane	226	C16H34	000544-76-3	87
3		Tetradecane	198	C14H30	000629-59-4	86
4		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	80
5		Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

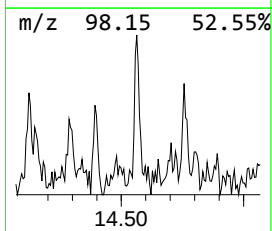
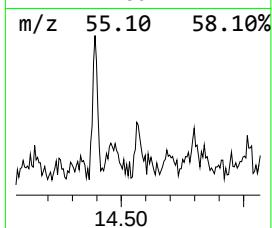
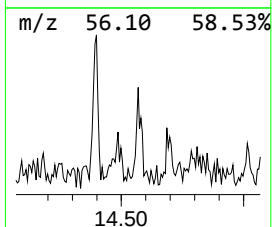
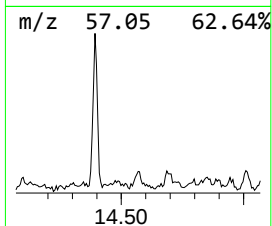
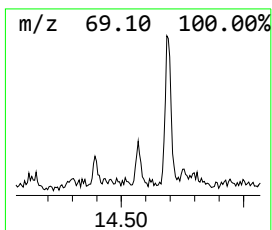
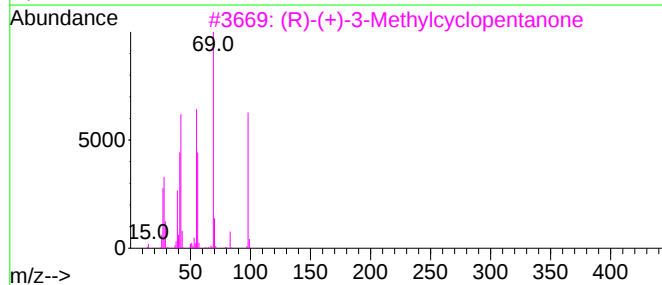
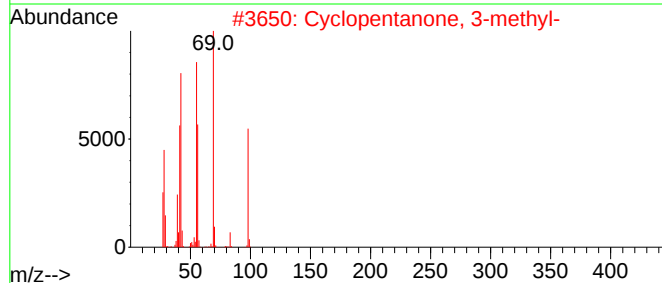
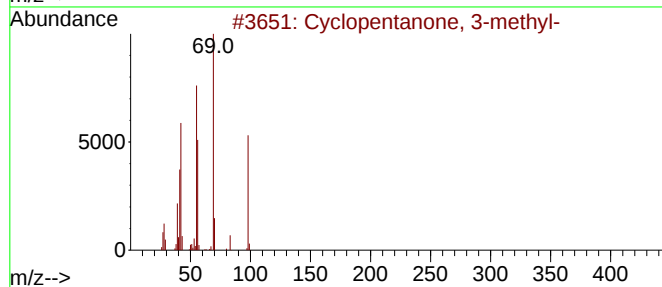
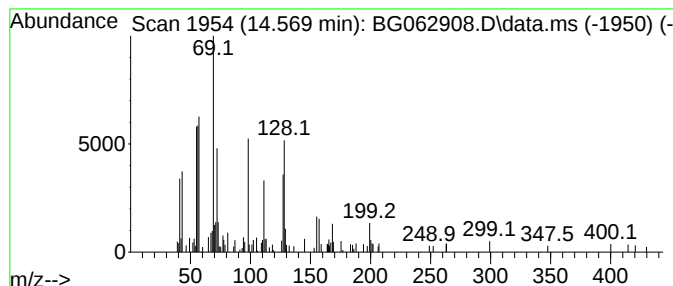
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 unknown-05 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.569	2.37 ng/ul	89512	Acenaphthene-d10			14.486
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopentanone, 3-methyl-		98	C6H10O	001757-42-2	43
2	Cyclopentanone, 3-methyl-		98	C6H10O	001757-42-2	43
3	(R)-(+)-3-Methylcyclopentanone		98	C6H10O	006672-30-6	38
4	Cyclohexane, 1,1,2-trimethyl-		126	C9H18	007094-26-0	38
5	Cyclobutanone, 2-(1,1-dimethylet...		126	C8H14O	004579-31-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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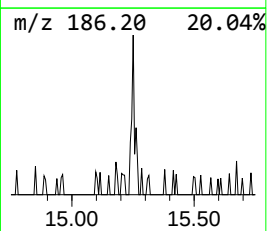
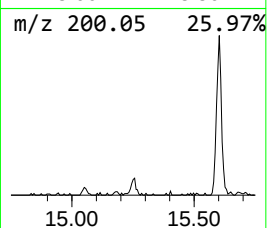
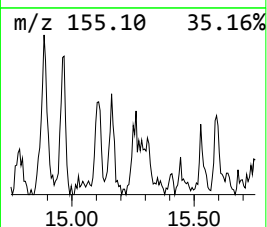
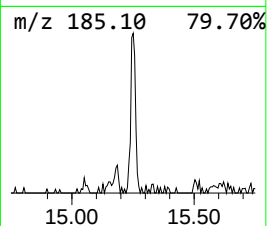
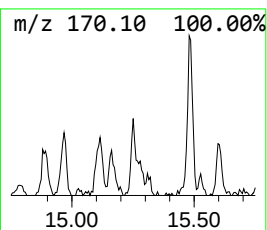
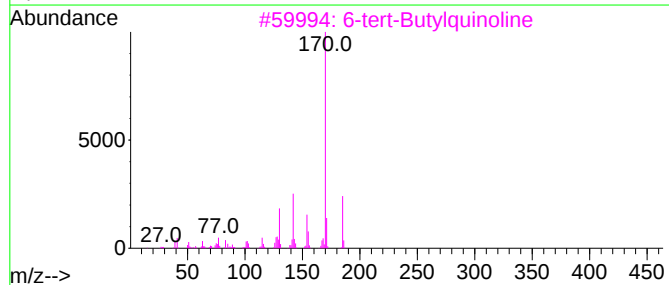
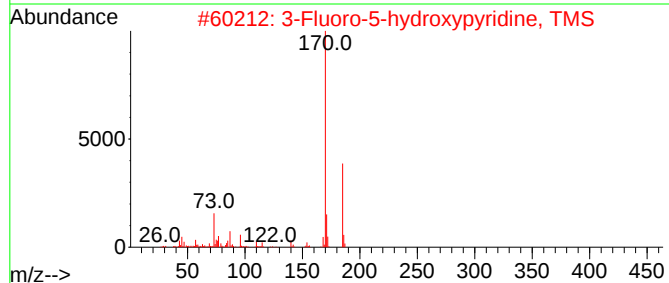
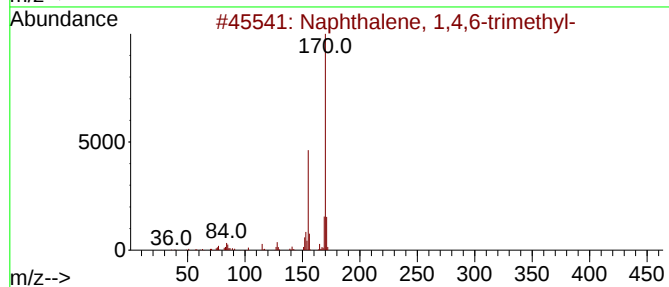
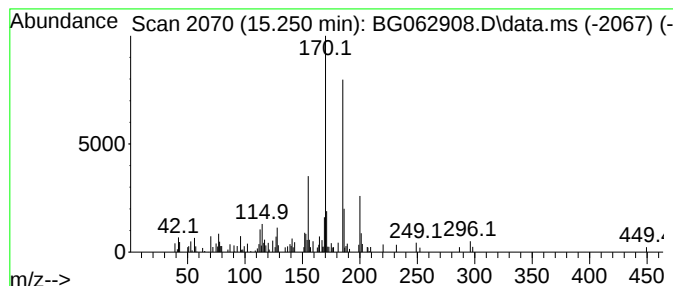
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Naphthalene, 1,4,6-trimethyl- Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.250	2.14 ng/ul	81084	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	83
2			3-Fluoro-5-hydroxypyridine, TMS	185	C8H12FNOSi	1000494-89-5	53
3			6-tert-Butylquinoline	185	C13H15N	068141-13-9	53
4			3,4-Dimethyl(1H)pyrrole, 2-[(3,4...	200	C13H16N2	065539-79-9	49
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

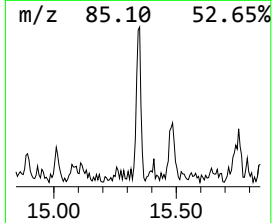
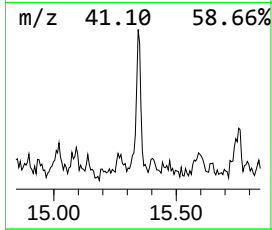
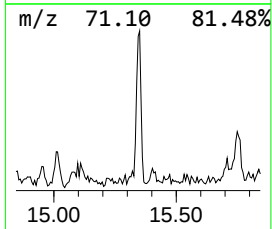
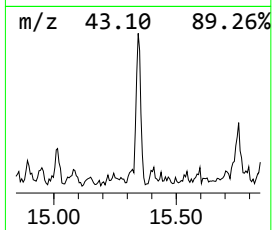
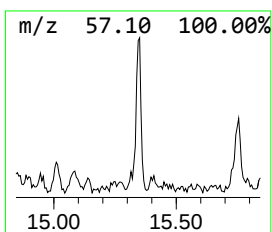
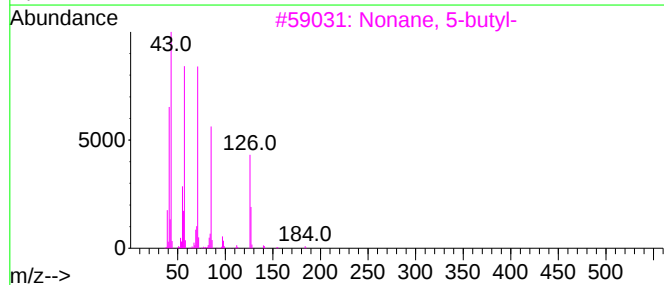
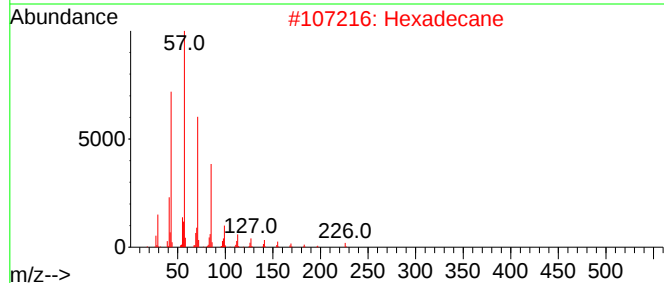
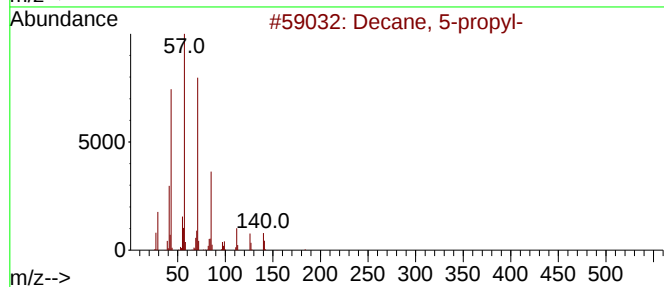
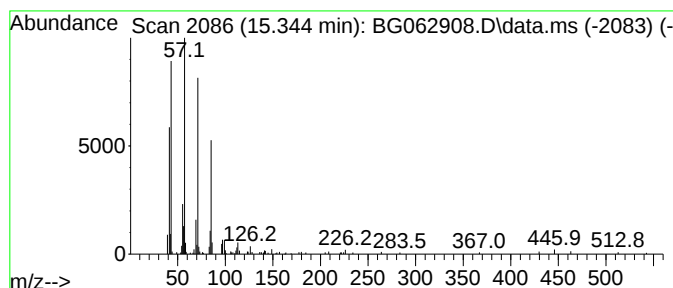
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.344	4.89 ng/ul	184885	Acenaphthene-d10	14.486	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5-propyl-	184	C13H28	017312-62-8	90
2	Hexadecane	226	C16H34	000544-76-3	87
3	Nonane, 5-butyl-	184	C13H28	017312-63-9	74
4	2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	64
5	Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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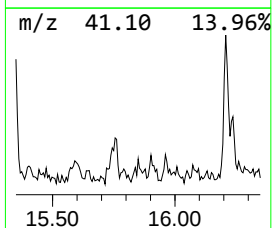
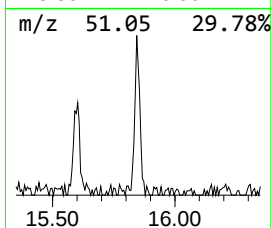
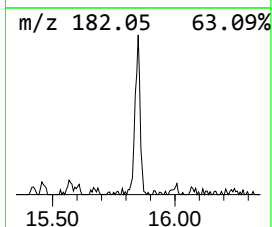
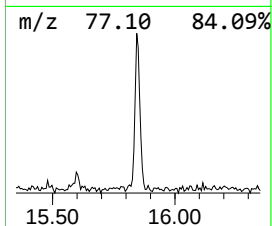
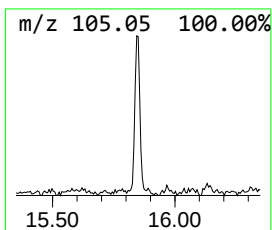
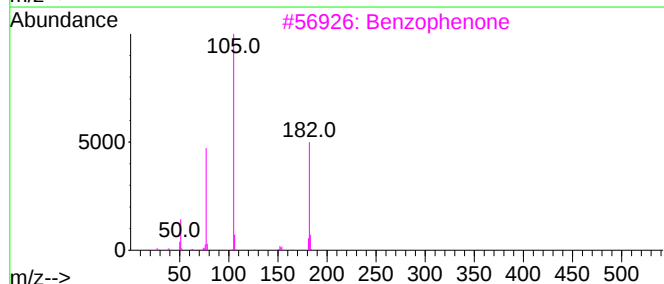
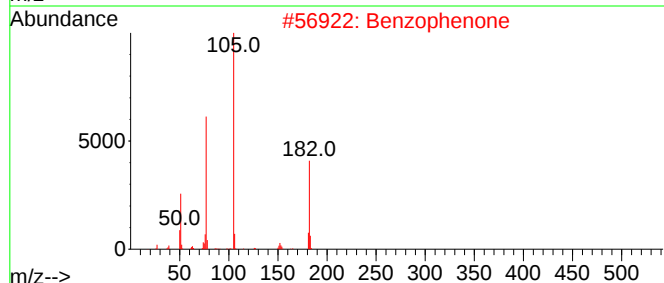
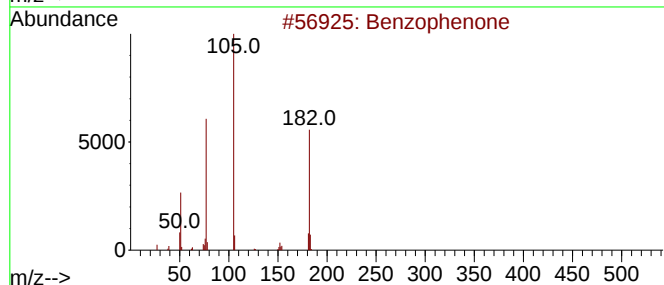
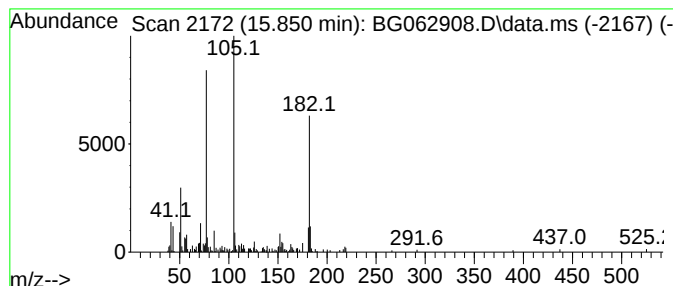
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Benzophenone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.850	4.02 ng/ul	151949	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	93
2			Benzophenone	182	C13H10O	000119-61-9	87
3			Benzophenone	182	C13H10O	000119-61-9	87
4			Benzophenone	182	C13H10O	000119-61-9	64
5			Benzoylformic acid	150	C8H6O3	000611-73-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

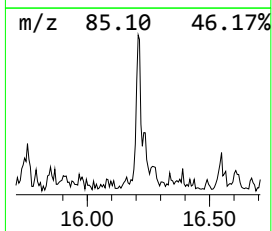
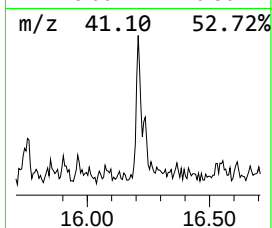
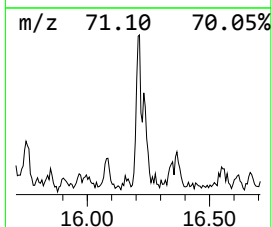
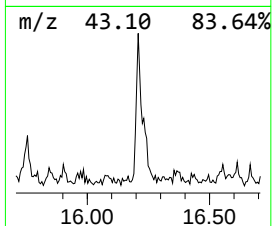
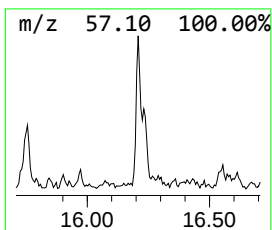
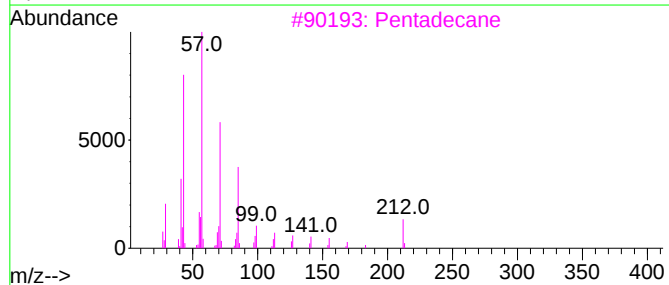
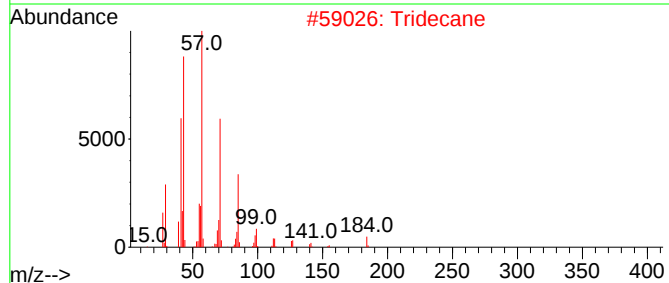
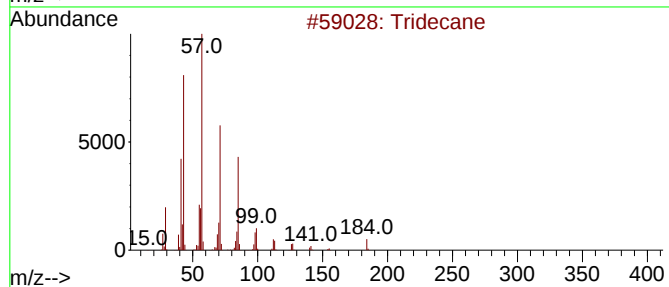
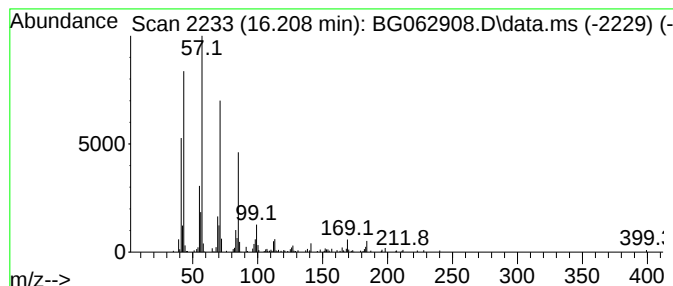
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.208	4.74 ng/ul	217593	Phenanthrene-d10	17.242

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	93
2		Tridecane	184	C13H28	000629-50-5	91
3		Pentadecane	212	C15H32	000629-62-9	87
4		Octacosane	394	C28H58	000630-02-4	83
5		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

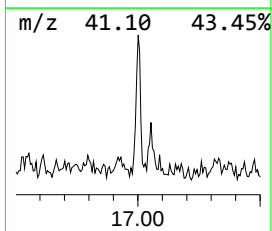
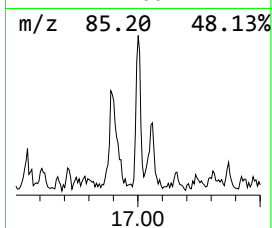
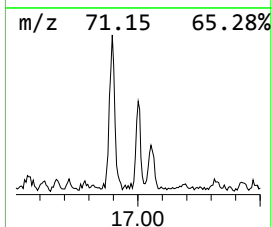
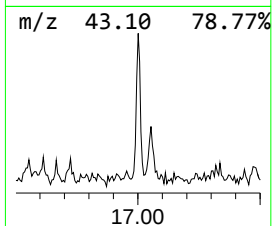
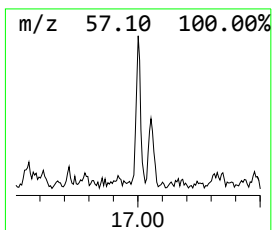
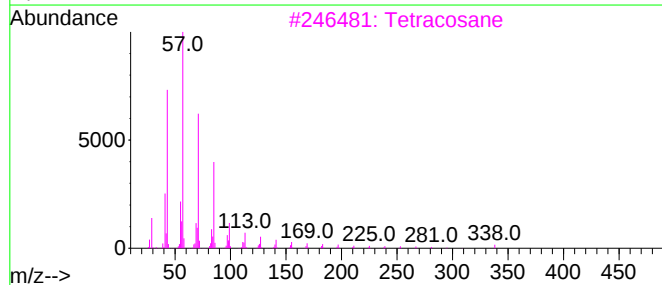
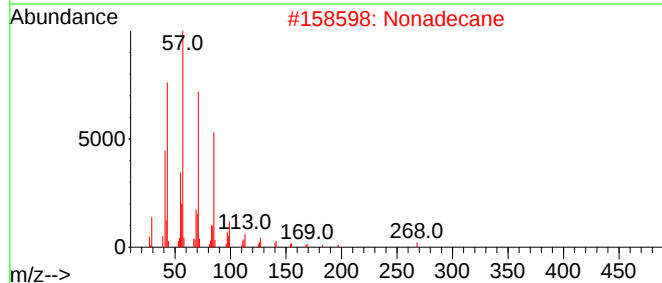
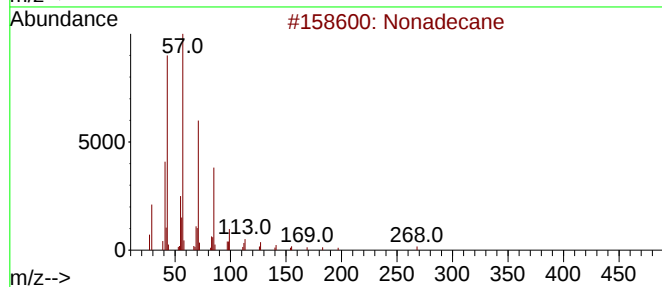
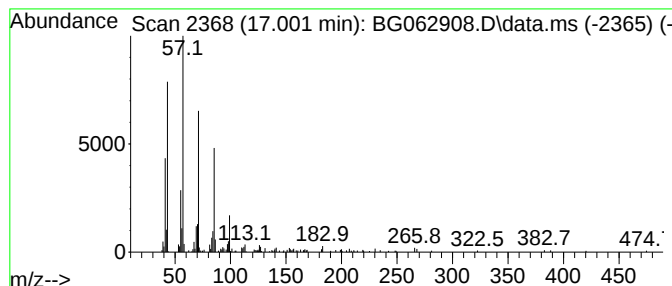
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.001	3.93 ng/ul	180099	Phenanthrene-d10	17.242

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	93
2		Nonadecane	268	C19H40	000629-92-5	91
3		Tetracosane	338	C24H50	000646-31-1	78
4		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	78
5		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

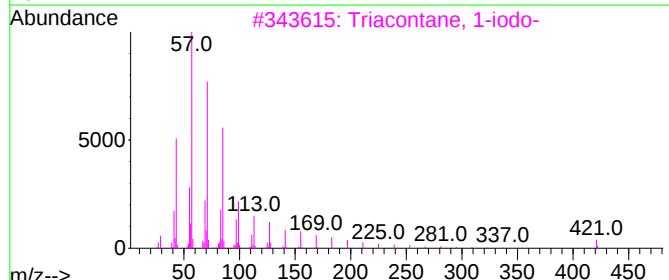
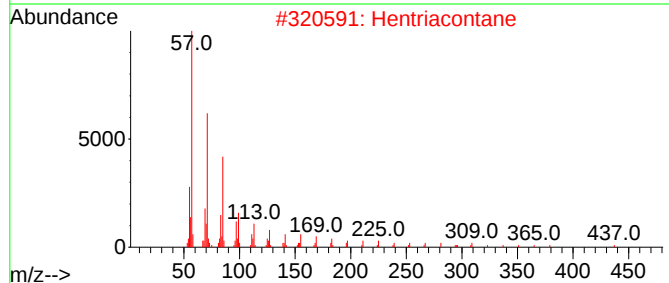
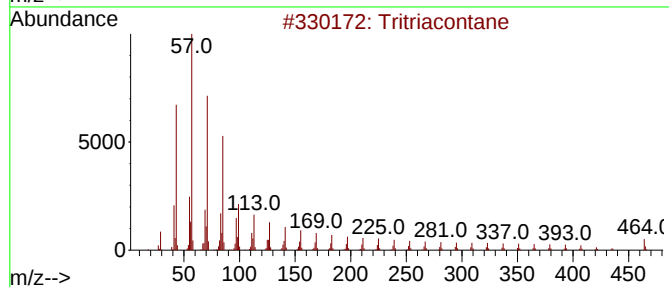
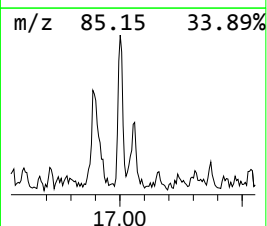
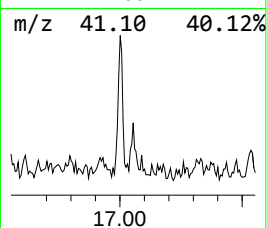
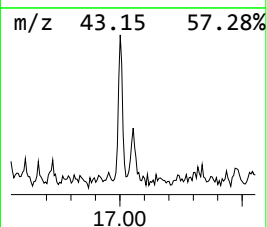
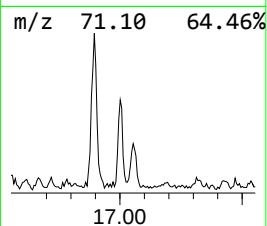
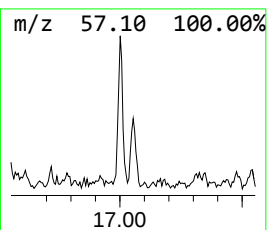
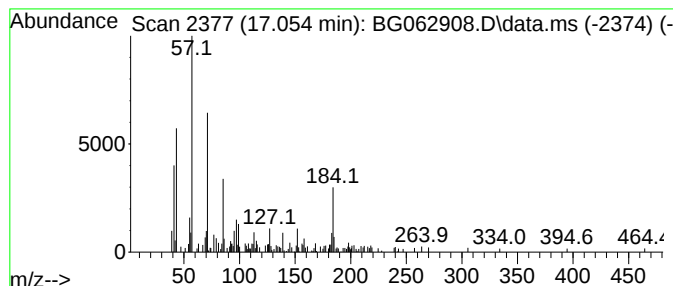
Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.054	2.22 ng/ul	101786	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tritriacontane		465	C33H68	000630-05-7	53
2	Hentriacontane		437	C31H64	000630-04-6	52
3	Triacontane, 1-iodo-		548	C30H61I	1000406-32-3	52
4	Hexacosane		366	C26H54	000630-01-3	52
5	Octacosane, 1-iodo-		520	C28H57I	1000406-32-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

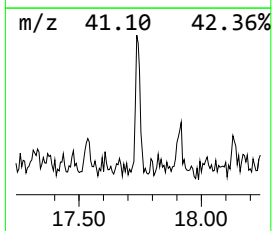
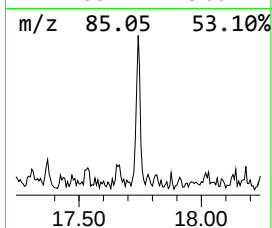
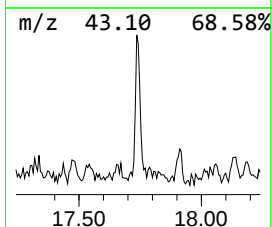
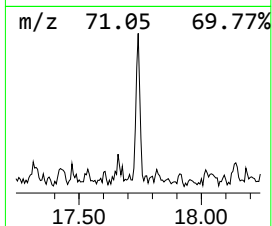
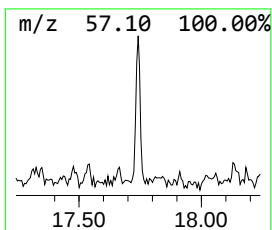
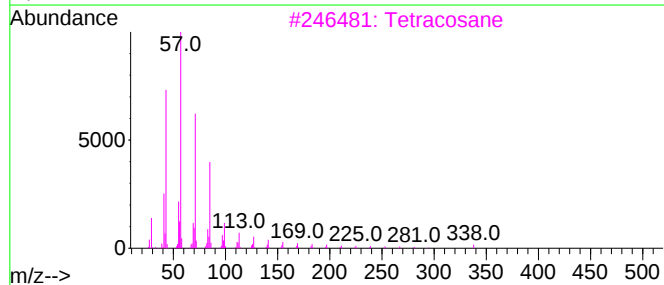
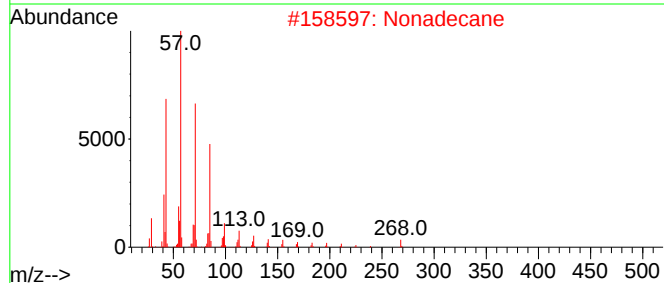
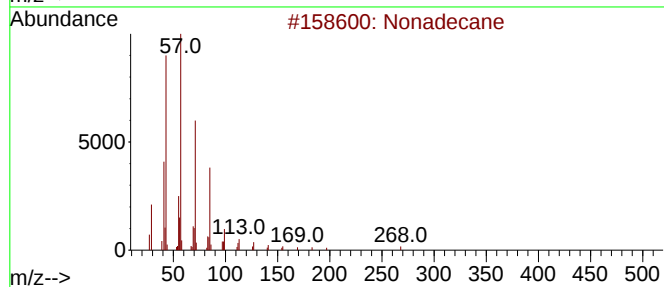
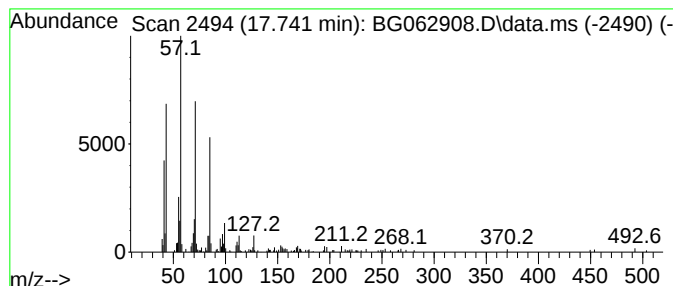
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.741	3.28 ng/ul	150304	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	92
2			Nonadecane	268	C19H40	000629-92-5	91
3			Tetracosane	338	C24H50	000646-31-1	83
4			Heneicosane	296	C21H44	000629-94-7	83
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

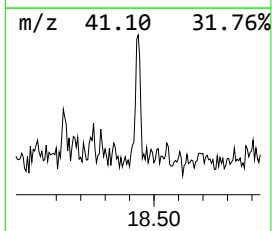
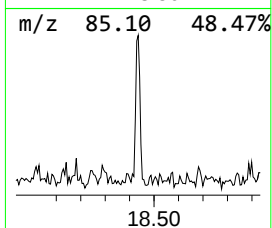
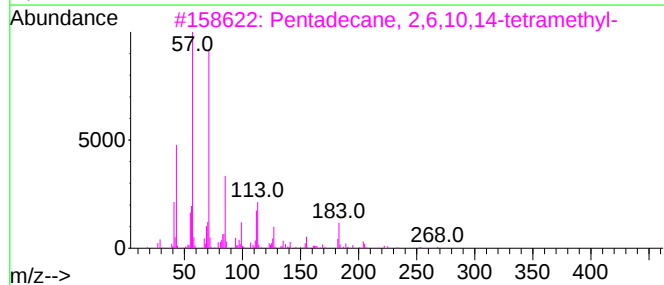
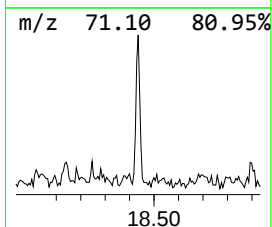
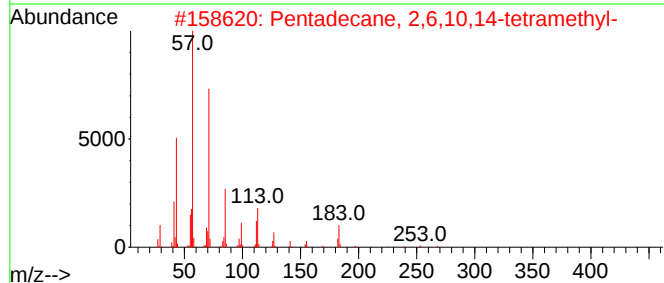
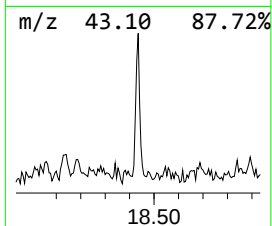
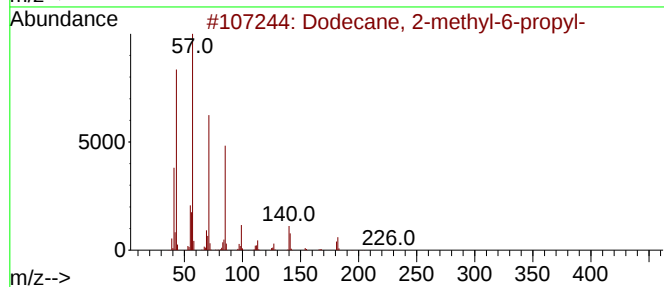
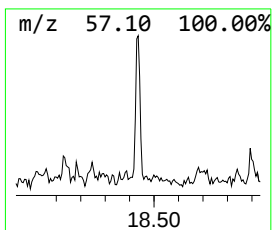
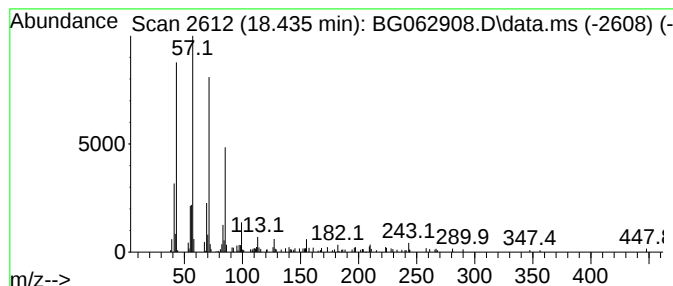
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.435	2.55 ng/ul	116780	Phenanthrene-d10			17.242
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
2	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	81
3	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	74
4	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	74
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

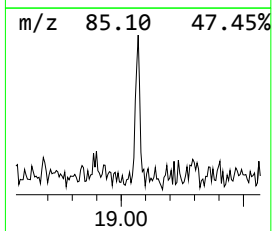
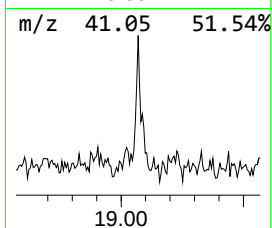
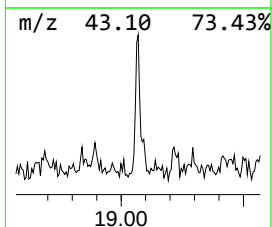
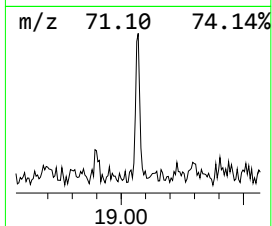
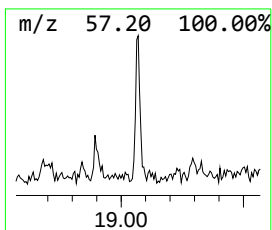
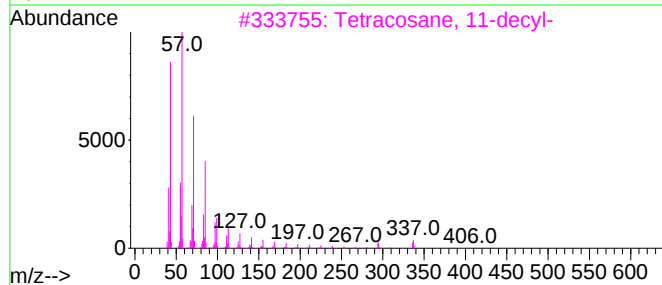
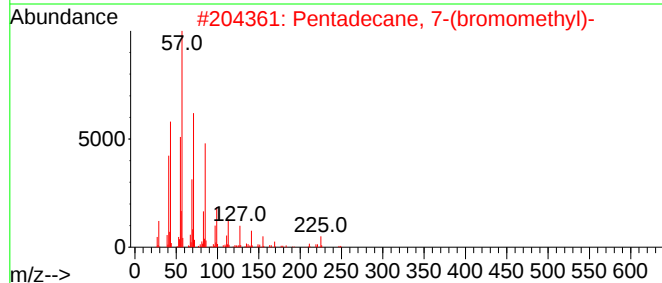
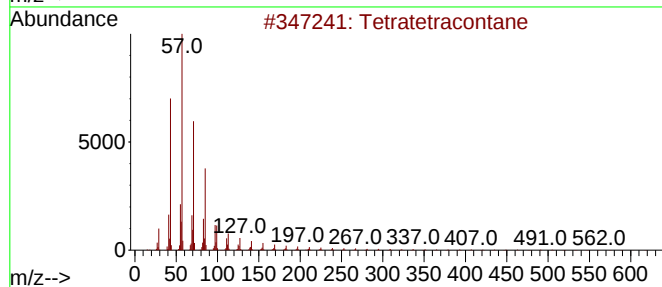
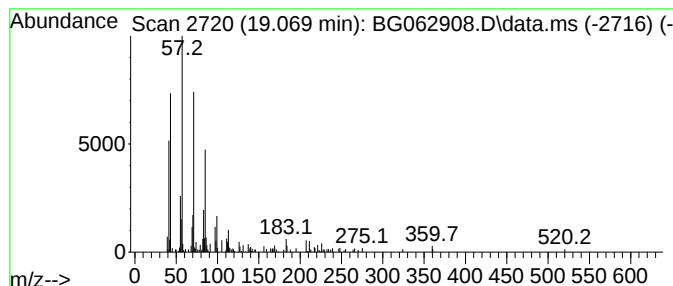
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.069	3.56 ng/ul	163456	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetratetracontane		619	C44H90	007098-22-8	90
2	Pentadecane, 7-(bromomethyl)-		304	C16H33Br	052997-43-0	90
3	Tetracosane, 11-decyl-		479	C34H70	055429-84-0	90
4	Trtriacontane, 2-methyl-		479	C34H70	066214-27-5	86
5	Heptacosane		380	C27H56	000593-49-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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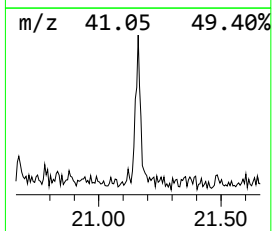
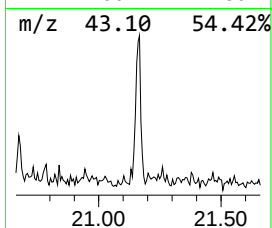
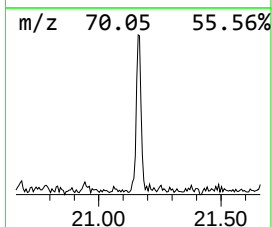
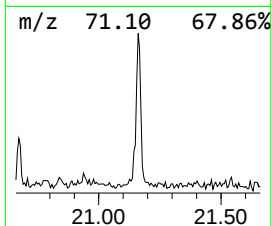
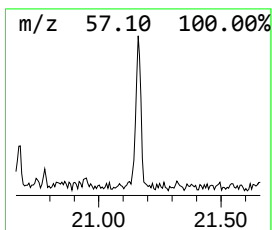
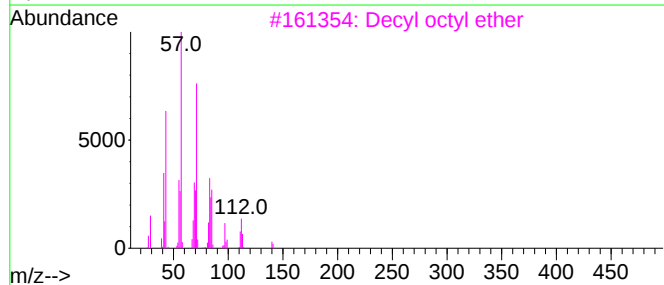
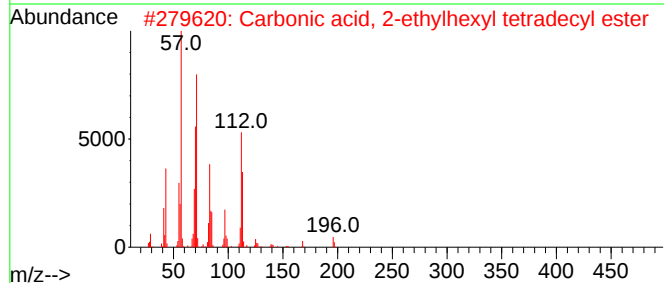
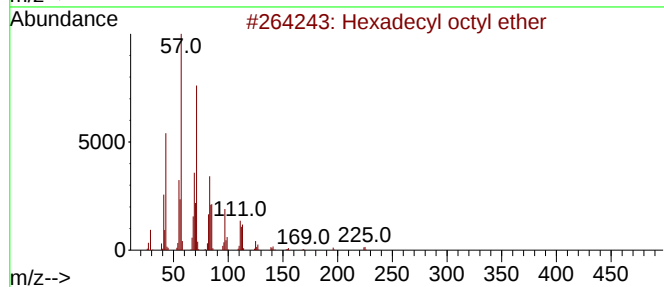
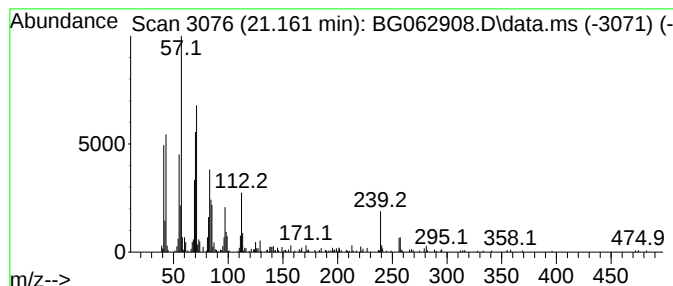
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 Hexadecyl octyl ether Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.161	6.28 ng/ul	313236	Chrysene-d12	21.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecyl octyl ether	354	C24H50O	1000406-38-6	74
2			Carbonic acid, 2-ethylhexyl tetr...	370	C23H46O3	1010383-14-1	64
3			Decyl octyl ether	270	C18H38O	1000406-38-3	64
4			Eicosyl octyl ether	410	C28H58O	1000406-38-8	58
5			Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062908.D
Acq On : 18 Sep 2024 14:29
Operator : JU/RC
Sample : P3878-12ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ0ME

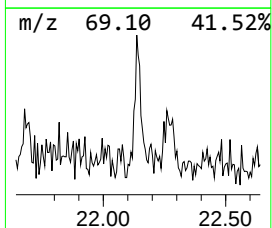
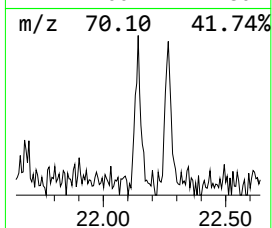
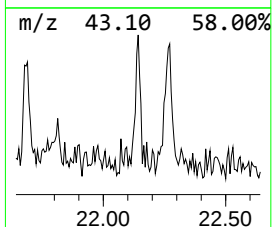
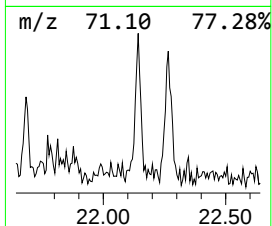
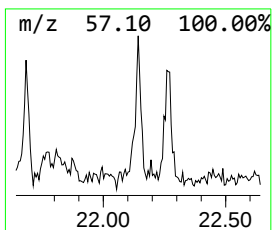
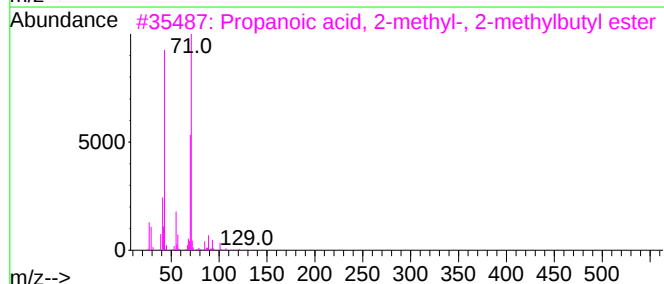
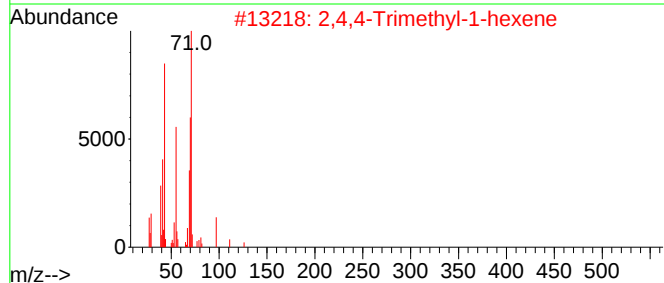
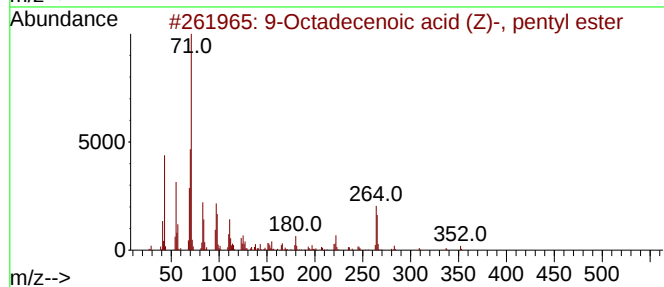
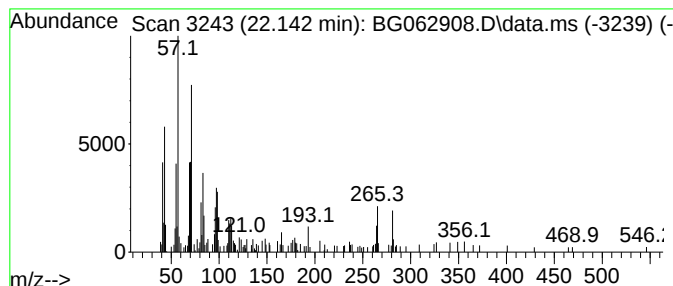
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 unknown-06 Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.142	2.22 ng/ul	110892	Chrysene-d12	21.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, pentyl...	352	C23H44O2	000142-57-4	47
2			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	43
3			Propanoic acid, 2-methyl-, 2-met...	158	C9H18O2	002445-69-4	38
4			n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	38
5			Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062908.D
 Acq On : 18 Sep 2024 14:29
 Operator : JU/RC
 Sample : P3878-12ME
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCVZ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.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...	5.891	6.0	ng/ul	110315	1	7.859	370673	20.0
Hexylene glycol	6.173	7.8	ng/ul	144703	1	7.859	370673	20.0
(DEL) Alkane: S...	6.361	2.8	ng/ul	52655	1	7.859	370673	20.0
(DEL) Alkane: S...	6.425	2.1	ng/ul	39135	1	7.859	370673	20.0
(DEL) Alkane: C...	6.502	3.0	ng/ul	55869	1	7.859	370673	20.0
unknown-01	6.766	2.0	ng/ul	37638	1	7.859	370673	20.0
(DEL) Alkane: S...	6.854	4.2	ng/ul	78028	1	7.859	370673	20.0
(DEL) Alkane: S...	6.913	4.3	ng/ul	79075	1	7.859	370673	20.0
Benzene, 1-ethy...	6.954	3.7	ng/ul	68876	1	7.859	370673	20.0
Mesitylene	7.083	2.5	ng/ul	45833	1	7.859	370673	20.0
Benzene, 1,2,4-...	7.248	2.3	ng/ul	42166	1	7.859	370673	20.0
2-Nonanone, 9-[...	7.289	2.4	ng/ul	44645	1	7.859	370673	20.0
(DEL) Alkane: S...	7.489	26.7	ng/ul	495105	1	7.859	370673	20.0
2-Propyl-1-pent...	7.918	11.0	ng/ul	204565	1	7.859	370673	20.0
Benzene, 1,2,3-...	7.976	3.0	ng/ul	55450	1	7.859	370673	20.0
unknown-02	8.147	2.2	ng/ul	40242	1	7.859	370673	20.0
Cyclohexanone, ...	8.282	3.8	ng/ul	69688	1	7.859	370673	20.0
(DEL) Alkane: S...	8.394	4.6	ng/ul	84810	1	7.859	370673	20.0
Methoxyacetic a...	8.452	2.2	ng/ul	40340	1	7.859	370673	20.0
1,3,8-p-Menthat...	8.499	4.5	ng/ul	84161	1	7.859	370673	20.0
(DEL) Alkane: S...	8.623	3.7	ng/ul	68129	1	7.859	370673	20.0
Benzene, 1-meth...	8.658	3.6	ng/ul	67490	1	7.859	370673	20.0
Benzene, 2-ethy...	8.811	2.6	ng/ul	48231	1	7.859	370673	20.0
Benzene, 1,2,4,...	8.858	3.1	ng/ul	57601	1	7.859	370673	20.0
Carbonic acid, ...	9.087	20.0	ng/ul	370695	1	7.859	370673	20.0
unknown-03	9.210	5.7	ng/ul	104863	1	7.859	370673	20.0
2,4-Dipropyl-5-...	11.020	5.8	ng/ul	221975	2	10.644	772289	20.0
(DEL) Alkane: S...	11.149	2.9	ng/ul	111453	2	10.644	772289	20.0
Cyclohexanol, 1...	11.672	2.6	ng/ul	99522	2	10.644	772289	20.0
Benzene, 1,3,5-...	11.743	2.3	ng/ul	86842	2	10.644	772289	20.0
1,3-Hexanediol,...	11.913	3.0	ng/ul	114182	2	10.644	772289	20.0
(DEL) Alkane: S...	12.072	4.5	ng/ul	174875	2	10.644	772289	20.0
unknown-04	12.107	2.9	ng/ul	111294	2	10.644	772289	20.0
(DEL) Alkane: S...	13.317	5.4	ng/ul	204750	3	14.486	756524	20.0
Naphthalene, 2,...	13.670	2.7	ng/ul	103606	3	14.486	756524	20.0
Naphthalene, 1,...	13.811	3.0	ng/ul	115283	3	14.486	756524	20.0
(DEL) Alkane: S...	14.392	6.3	ng/ul	238515	3	14.486	756524	20.0
unknown-05	14.569	2.4	ng/ul	89512	3	14.486	756524	20.0
Naphthalene, 1,...	15.250	2.1	ng/ul	81084	3	14.486	756524	20.0
(DEL) Alkane: S...	15.344	4.9	ng/ul	184885	3	14.486	756524	20.0
Benzophenone	15.850	4.0	ng/ul	151949	3	14.486	756524	20.0
(DEL) Alkane: S...	16.208	4.7	ng/ul	217593	4	17.242	917196	20.0
(DEL) Alkane: S...	17.001	3.9	ng/ul	180099	4	17.242	917196	20.0
(DEL) Alkane: S...	17.054	2.2	ng/ul	101786	4	17.242	917196	20.0
(DEL) Alkane: S...	17.741	3.3	ng/ul	150304	4	17.242	917196	20.0
(DEL) Alkane: S...	18.435	2.5	ng/ul	116780	4	17.242	917196	20.0
(DEL) Alkane: S...	19.069	3.6	ng/ul	163456	4	17.242	917196	20.0
Hexadecyl octyl...	21.161	6.3	ng/ul	313236	5	21.490	997984	20.0
unknown-06	22.142	2.2	ng/ul	110892	5	21.490	997984	20.0