

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062949.D
 Acq On : 19 Sep 2024 18:49
 Operator : JU/RC
 Sample : P3878-15MEDL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC04MEDL

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: BG062949.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.954	145	147	148	rBV3	5506	3177	0.33%	0.047%
2	4.477	235	236	239	rBV	5496	4238	0.43%	0.062%
3	4.712	274	276	277	rBV	3563	2721	0.28%	0.040%
4	5.035	328	331	332	rBV2	3345	3095	0.32%	0.046%
5	5.088	338	340	344	rBV4	2840	2929	0.30%	0.043%
6	5.393	390	392	395	rVB2	2944	2539	0.26%	0.037%
7	5.428	395	398	399	rBV2	2683	3034	0.31%	0.045%
8	5.663	436	438	442	rVB2	3613	3344	0.34%	0.049%
9	5.810	460	463	464	rBV2	3677	3623	0.37%	0.053%
10	5.887	469	476	483	rBV5	18008	36966	3.79%	0.544%
11	6.251	536	538	541	rBV4	5203	6178	0.63%	0.091%
12	6.357	551	556	558	rBV5	7745	8865	0.91%	0.130%
13	6.433	564	569	572	rBV5	8302	14400	1.48%	0.212%
14	6.474	574	576	579	rBV2	3113	4217	0.43%	0.062%
15	6.603	594	598	599	rBV2	3957	3674	0.38%	0.054%
16	6.656	604	607	608	rBV2	3781	4209	0.43%	0.062%
17	6.692	612	613	615	rVB	5536	2620	0.27%	0.039%
18	6.739	619	621	622	rBV2	6095	4397	0.45%	0.065%
19	6.856	633	641	645	rBV5	14142	33584	3.44%	0.494%
20	6.909	648	650	654	rVV2	13203	15879	1.63%	0.234%
21	7.027	664	670	674	rBV5	34040	66182	6.78%	0.974%
22	7.085	677	680	687	rVB2	12655	19832	2.03%	0.292%
23	7.185	694	697	702	rVB2	12898	19435	1.99%	0.286%
24	7.244	702	707	711	rBV5	15005	22033	2.26%	0.324%
25	7.485	743	748	756	rVB3	91457	161385	16.53%	2.374%
26	7.855	804	811	817	rBV2	156268	277882	28.46%	4.088%
27	7.920	819	822	828	rVV3	23596	37974	3.89%	0.559%
28	7.967	828	830	839	rVB6	12059	21815	2.23%	0.321%
29	8.113	853	855	856	rBV2	4723	2928	0.30%	0.043%
30	8.202	868	870	872	rBV2	4263	4088	0.42%	0.060%
31	8.237	874	876	878	rVB2	5365	3189	0.33%	0.047%
32	8.284	878	884	888	rBV3	13852	27827	2.85%	0.409%
33	8.348	892	895	896	rBV3	6523	6184	0.63%	0.091%
34	8.448	909	912	915	rVB4	11517	14134	1.45%	0.208%
35	8.495	916	920	921	rBV3	17663	21636	2.22%	0.318%
36	8.619	937	941	945	rBV4	16463	25511	2.61%	0.375%

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Integration Parameters: LSCINT.P

Integrator: RTE

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37	8.813	969	974	977	rBV2	10595	19410	1.99%	0.286%
38	8.860	979	982	985	rVB4	10730	12120	1.24%	0.178%
39	8.924	991	993	994	rBV2	4469	2829	0.29%	0.042%
40	9.083	1016	1020	1026	rVB2	78109	114848	11.76%	1.689%
41	9.206	1032	1041	1046	rBV9	10691	31982	3.28%	0.470%
42	9.729	1123	1130	1135	rBV7	22452	45426	4.65%	0.668%
43	10.000	1173	1176	1180	rVB5	7936	11003	1.13%	0.162%
44	10.035	1180	1182	1184	rBV3	4072	4258	0.44%	0.063%
45	10.334	1229	1233	1239	rBV6	10823	17851	1.83%	0.263%
46	10.411	1243	1246	1248	rBV2	2758	2993	0.31%	0.044%
47	10.475	1252	1257	1259	rBV3	3391	4529	0.46%	0.067%
48	10.646	1276	1286	1292	rBV	209953	426160	43.65%	6.269%
49	10.769	1302	1307	1308	rBV4	18301	22267	2.28%	0.328%
50	10.916	1328	1332	1334	rBV3	3857	4881	0.50%	0.072%
51	11.022	1345	1350	1356	rVB3	36055	59863	6.13%	0.881%
52	11.151	1366	1372	1377	rBV4	15163	24944	2.56%	0.367%
53	11.263	1388	1391	1394	rBV3	4502	5949	0.61%	0.088%
54	11.363	1401	1408	1410	rBV4	6495	14172	1.45%	0.208%
55	11.551	1436	1440	1441	rBV3	7384	9551	0.98%	0.140%
56	11.674	1457	1461	1466	rBV2	16644	24806	2.54%	0.365%
57	11.909	1497	1501	1505	rBV2	14313	21465	2.20%	0.316%
58	11.980	1511	1513	1517	rVB3	4408	3545	0.36%	0.052%
59	12.074	1521	1529	1532	rBV2	25391	48834	5.00%	0.718%
60	12.314	1565	1570	1575	rVB3	38821	62261	6.38%	0.916%
61	12.520	1603	1605	1612	rVB4	8157	12276	1.26%	0.181%
62	12.643	1624	1626	1629	rBV3	3805	4315	0.44%	0.063%
63	12.667	1629	1630	1631	rBV3	4312	2705	0.28%	0.040%
64	12.978	1679	1683	1685	rBV4	4842	6790	0.70%	0.100%
65	13.025	1689	1691	1694	rBV3	10147	10491	1.07%	0.154%
66	13.313	1735	1740	1749	rBV2	32660	59507	6.10%	0.875%
67	13.372	1749	1750	1753	rVB	7086	4245	0.43%	0.062%
68	13.407	1753	1756	1759	rBV3	5152	6676	0.68%	0.098%
69	13.807	1819	1824	1829	rBV5	16223	32055	3.28%	0.472%
70	13.889	1836	1838	1845	rVB2	39081	53293	5.46%	0.784%
71	14.394	1920	1924	1928	rVB4	38938	51380	5.26%	0.756%
72	14.488	1934	1940	1946	rVB	334987	518102	53.07%	7.621%
73	14.841	1996	2000	2002	rBV5	7125	9852	1.01%	0.145%
74	15.011	2026	2029	2037	rVB8	9582	20005	2.05%	0.294%
75	15.082	2039	2041	2042	rBV2	7620	6107	0.63%	0.090%
76	15.117	2044	2047	2052	rVB5	44574	66863	6.85%	0.984%

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 Title : SVOA CALIBRATION

77	15.346	2083	2086	2091	rVB2	39878	51835	5.31%	0.763%
78	15.481	2105	2109	2115	rVB2	52982	80028	8.20%	1.177%
79	16.210	2229	2233	2235	rBV	47446	59207	6.06%	0.871%
80	16.891	2343	2349	2363	rBV	628160	976271	100.00%	14.361%
81	17.003	2365	2368	2373	rVB3	32314	43520	4.46%	0.640%
82	17.238	2402	2408	2417	rVB	445513	679394	69.59%	9.994%
83	17.338	2421	2425	2429	rVB2	62188	88781	9.09%	1.306%
84	17.738	2490	2493	2498	rVB2	33553	43224	4.43%	0.636%
85	18.437	2608	2612	2615	rVB3	28642	33170	3.40%	0.488%
86	19.065	2717	2719	2726	rVB6	23351	41235	4.22%	0.607%
87	19.629	2811	2815	2822	rVB	85680	138931	14.23%	2.044%
88	20.675	2991	2993	2995	rBV3	16285	14554	1.49%	0.214%
89	21.163	3073	3076	3082	rVB4	47387	60117	6.16%	0.884%
90	21.486	3126	3131	3141	rVB	488193	809941	82.96%	11.914%
91	24.530	3640	3649	3663	rBV	338600	921404	94.38%	13.554%

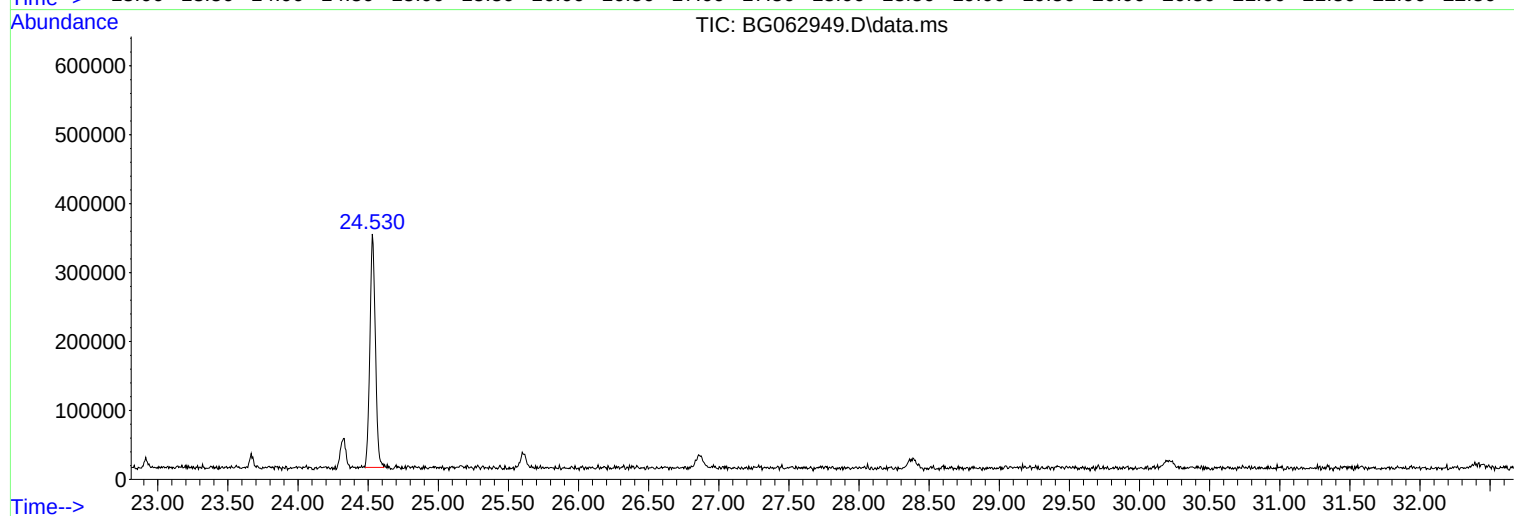
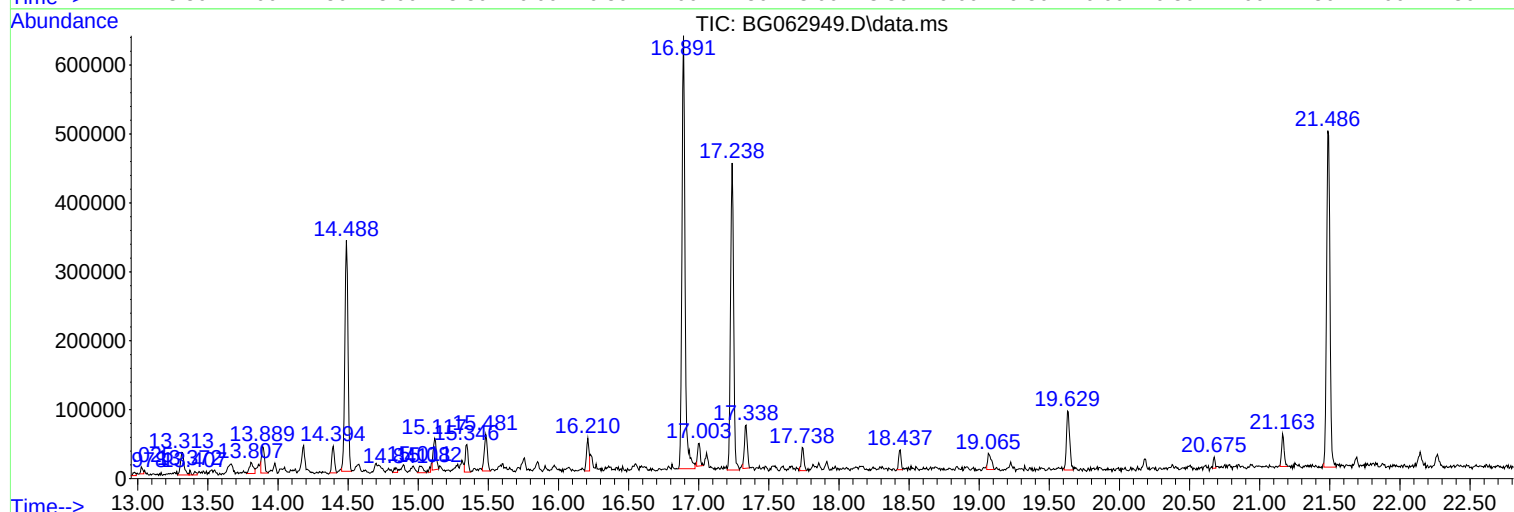
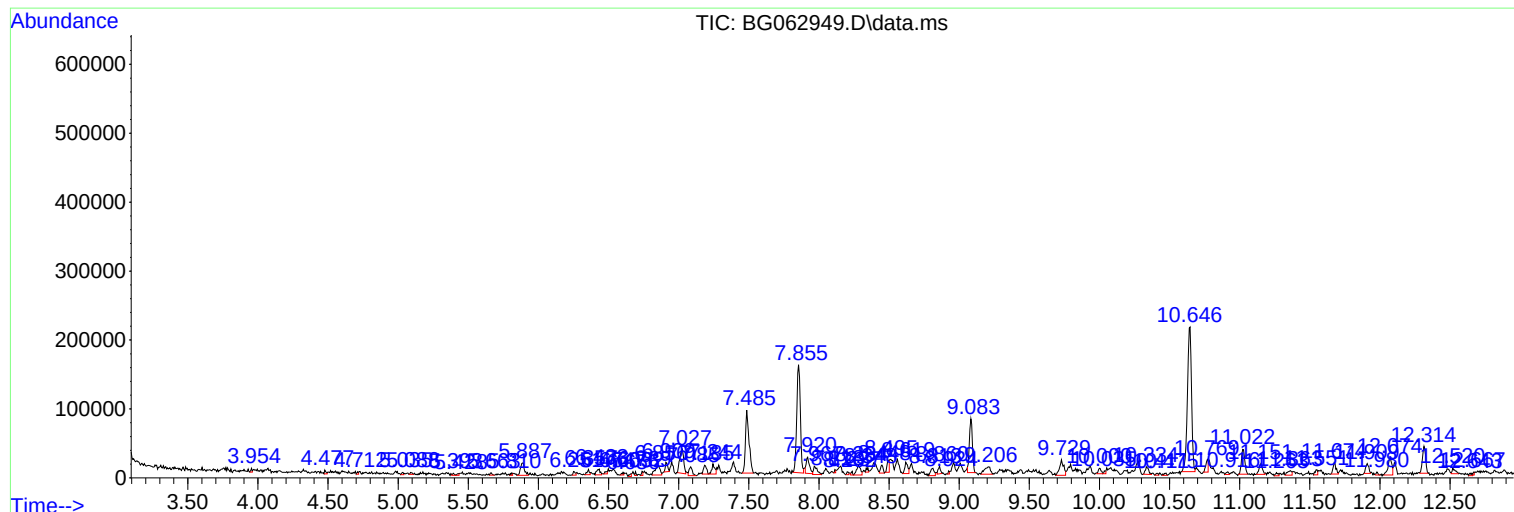
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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



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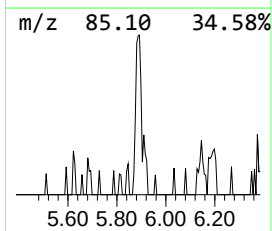
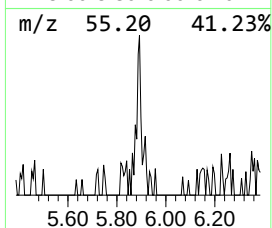
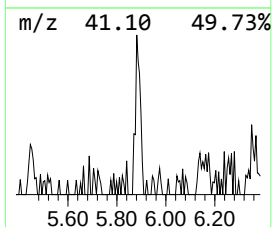
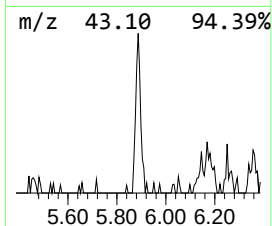
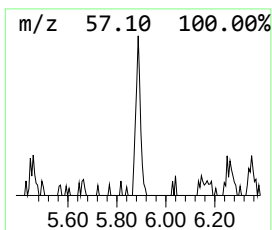
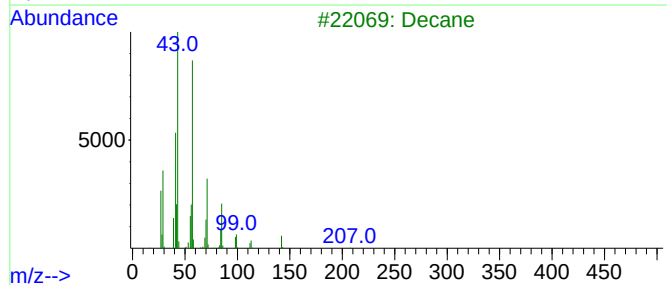
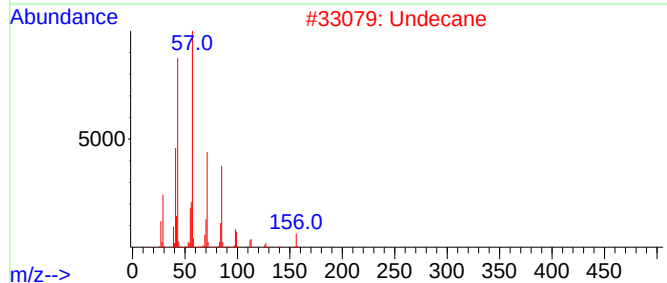
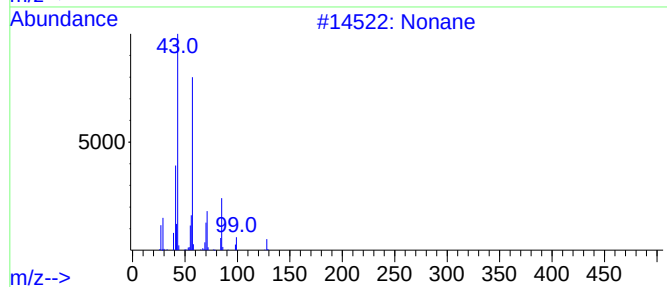
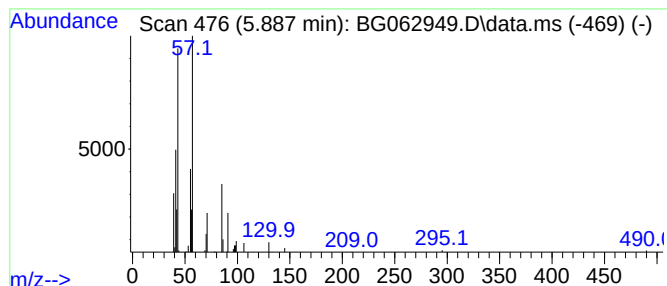
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.887	2.66 ng/ul	36966	1,4-Dichlorobenzene-d4	7.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	53
2			Undecane	156	C11H24	001120-21-4	50
3			Decane	142	C10H22	000124-18-5	40
4			Octane	114	C8H18	000111-65-9	40
5			Tridecane	184	C13H28	000629-50-5	40



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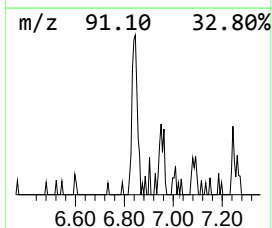
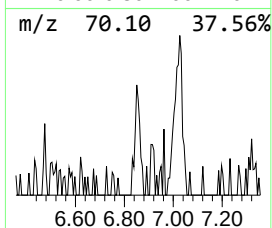
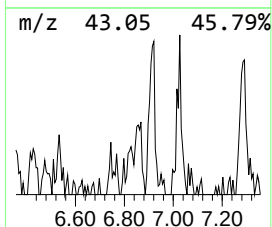
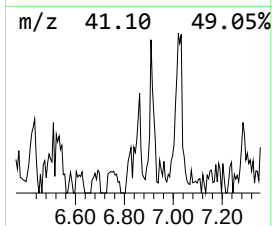
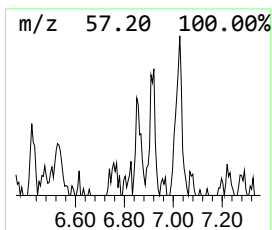
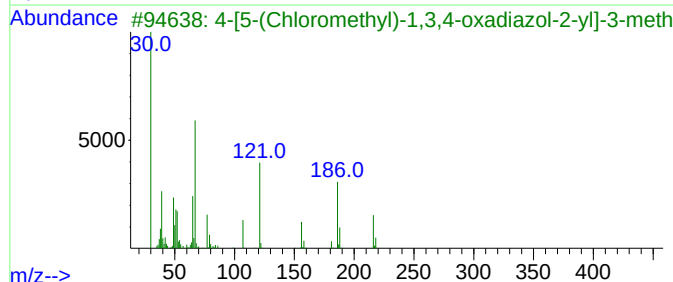
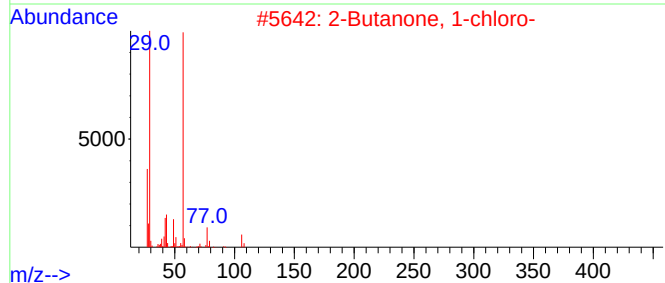
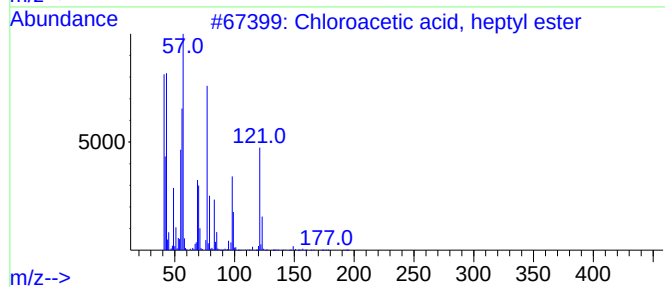
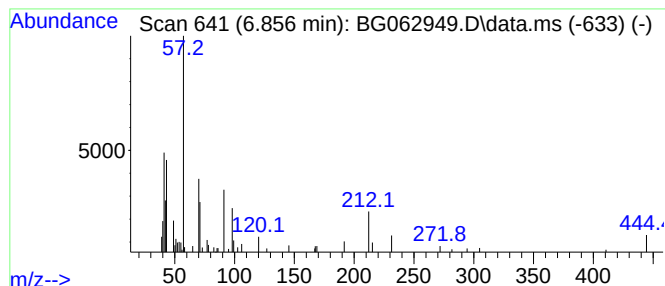
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.856	2.42 ng/ul	33584	1,4-Dichlorobenzene-d4	7.861

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Chloroacetic acid, heptyl ester	192	C9H17ClO2	034589-22-5	28
2		2-Butanone, 1-chloro-	106	C4H7ClO	000616-27-3	25
3		4-[5-(Chloromethyl)-1,3,4-oxadia...	216	C6H5ClN4O3	1000443-67-4	17
4		N-(2,2-Dichloro-1-hydroxy-ethyl)...	213	C7H13Cl2NO2	058956-78-8	9
5		.beta.-Chlordene	336	C10H6Cl6	056534-03-3	4



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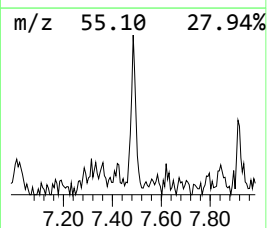
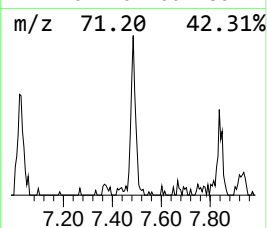
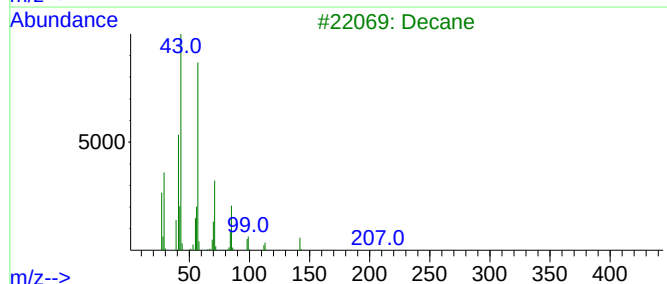
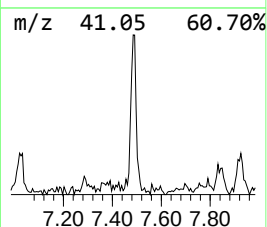
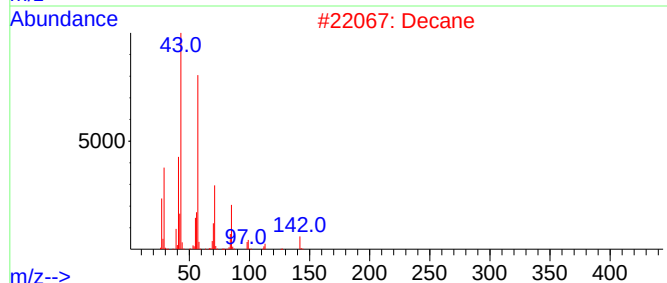
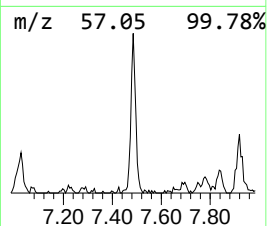
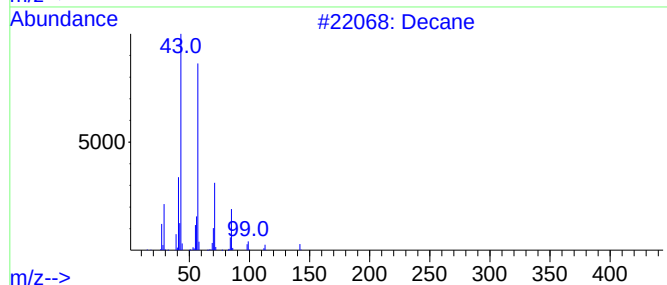
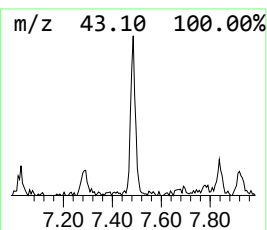
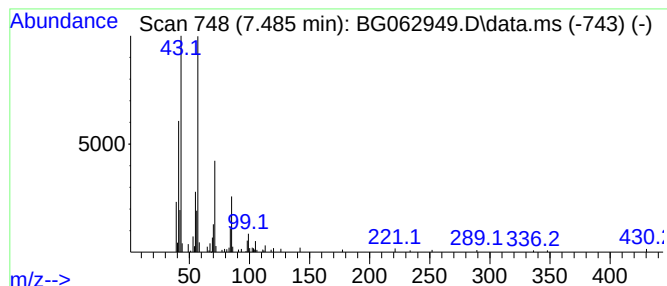
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.485	11.62 ng/ul	161385	1,4-Dichlorobenzene-d4	7.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane			142	C10H22	000124-18-5	87
2	Decane			142	C10H22	000124-18-5	86
3	Decane			142	C10H22	000124-18-5	78
4	Undecane			156	C11H24	001120-21-4	78
5	Tridecane			184	C13H28	000629-50-5	72



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Misc :
ALS Vial : 12 Sample Multiplier: 1

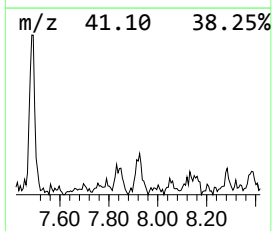
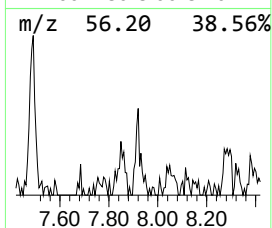
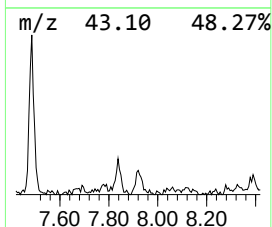
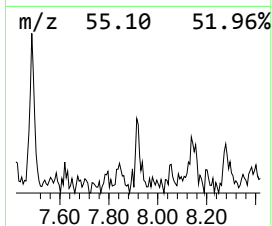
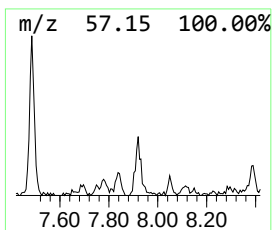
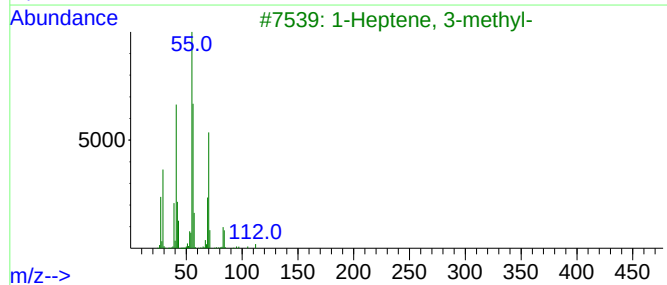
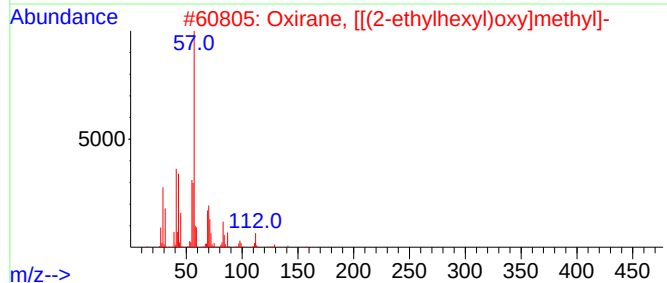
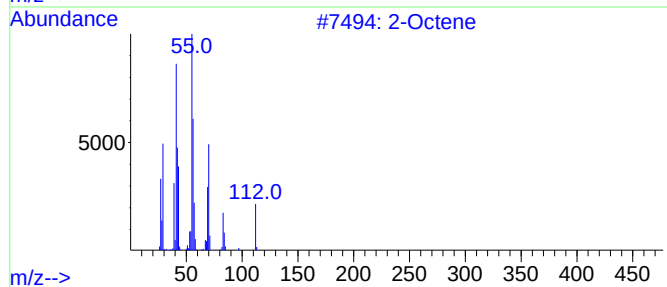
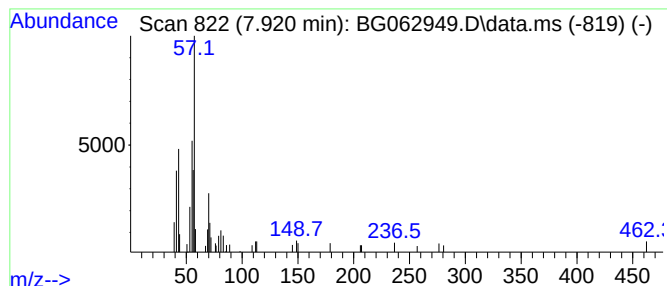
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BNA_G
ClientSampleId :
DDC04MEDL

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 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.920	2.73 ng/ul	37974	1,4-Dichlorobenzene-d4	7.861	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Octene	112	C8H16	000111-67-1	43
2	Oxirane, [[(2-ethylhexyl)oxy]met...	186	C11H22O2	002461-15-6	38
3	1-Heptene, 3-methyl-	112	C8H16	004810-09-7	32
4	6-Methyl-2-heptanol, trifluoroac...	226	C10H17F3O2	1010365-05-8	32
5	Isobutyl pentyl carbonate	188	C10H20O3	178442-32-5	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

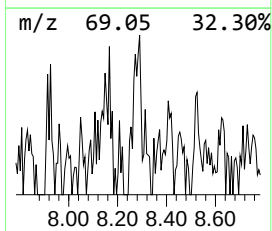
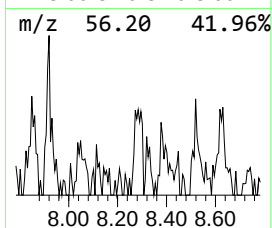
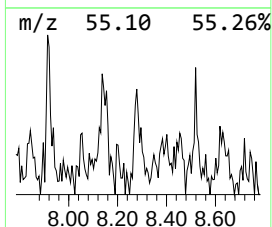
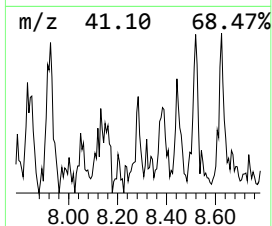
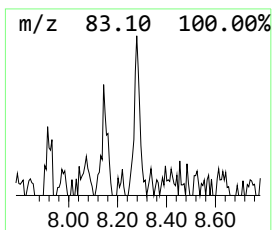
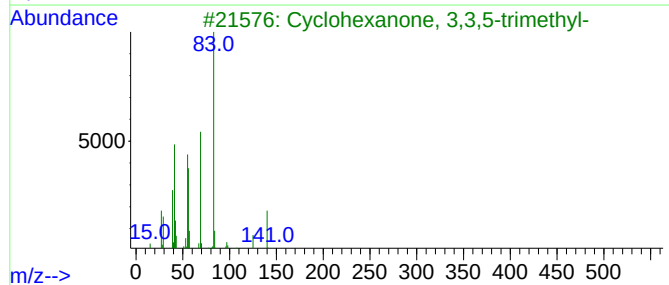
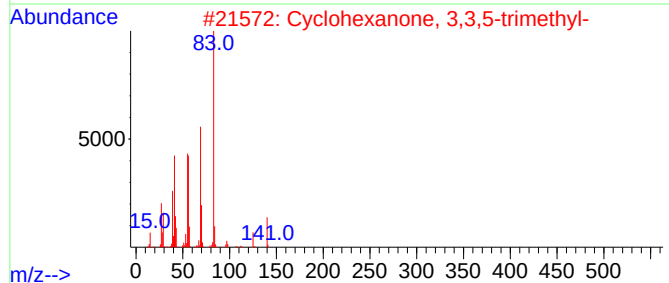
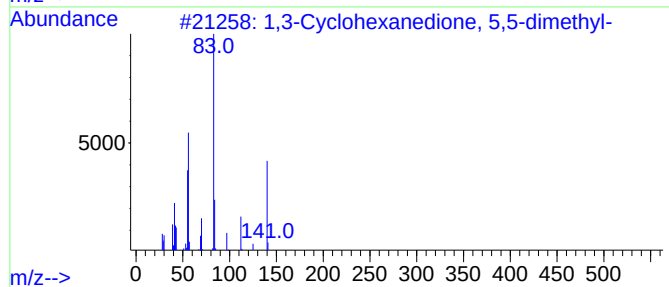
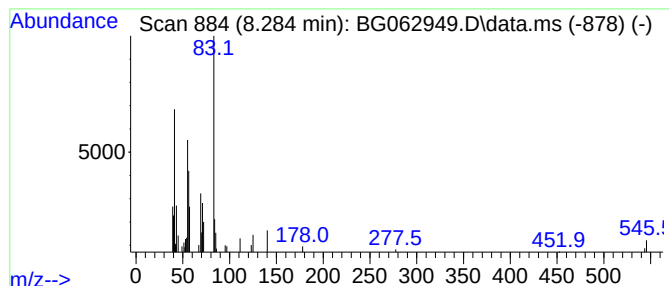
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,3-Cyclohexanedione, 5,5-d... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.284	2.00 ng/ul	27827	1,4-Dichlorobenzene-d4	7.861	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	50
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	47
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	47
4	Propanenitrile, 2-(dimethylamino)-	98	C5H10N2	005350-67-4	42
5	2,4,6-Trimethyl-3-heptene	140	C10H20	1000113-56-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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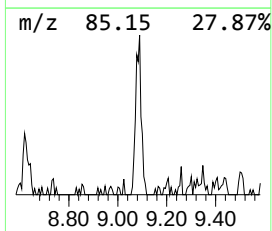
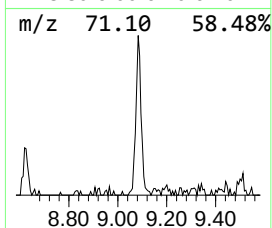
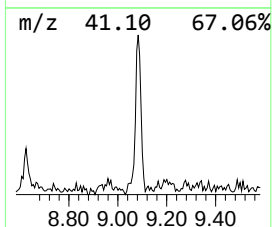
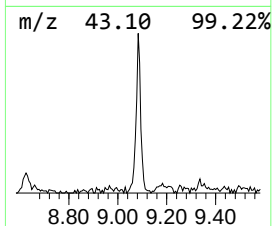
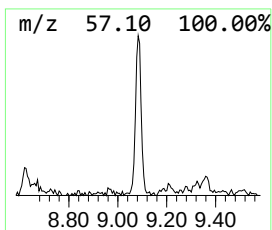
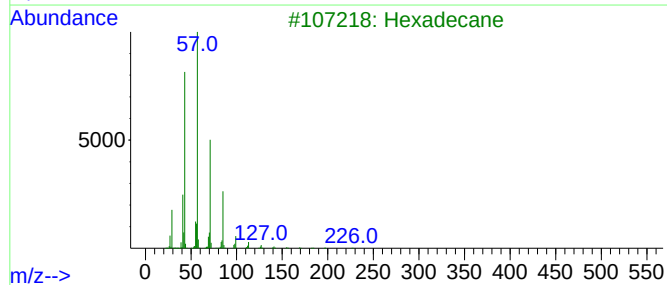
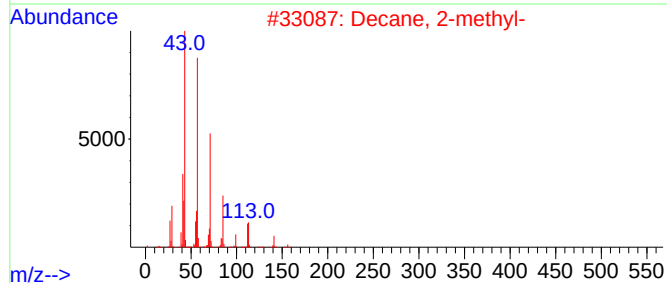
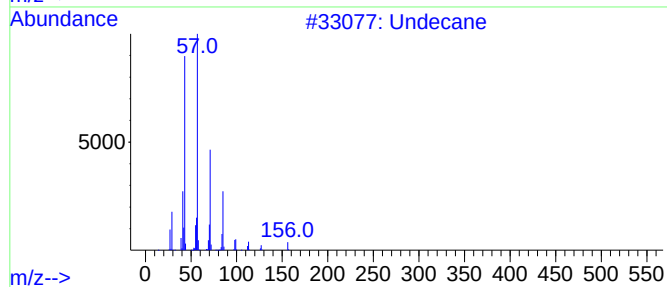
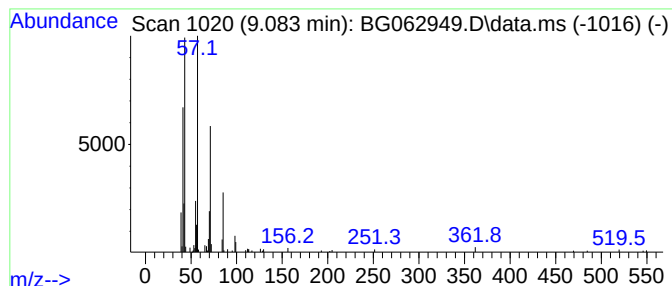
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.083	8.27 ng/ul	114848	1,4-Dichlorobenzene-d4	7.861

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	68
2	Decane, 2-methyl-	156	C11H24	006975-98-0	58
3	Hexadecane	226	C16H34	000544-76-3	58
4	Tridecane	184	C13H28	000629-50-5	58
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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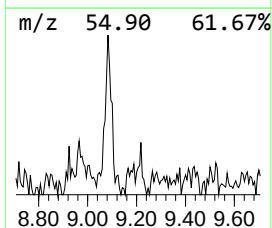
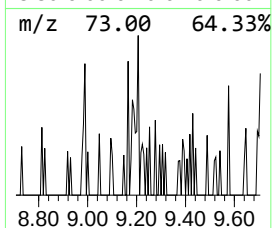
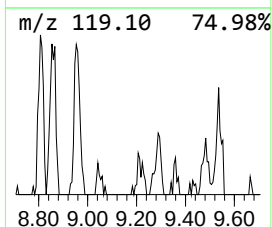
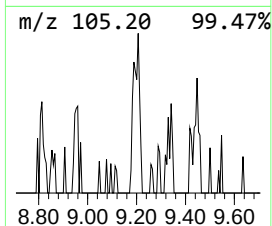
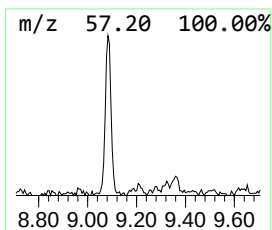
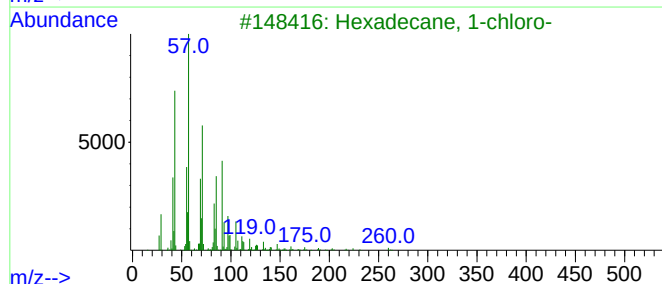
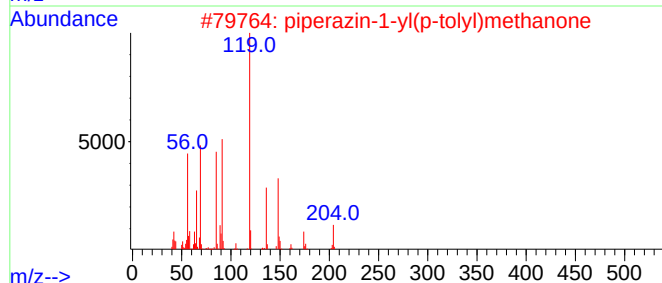
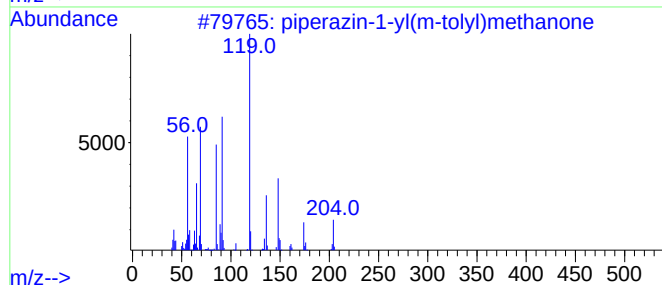
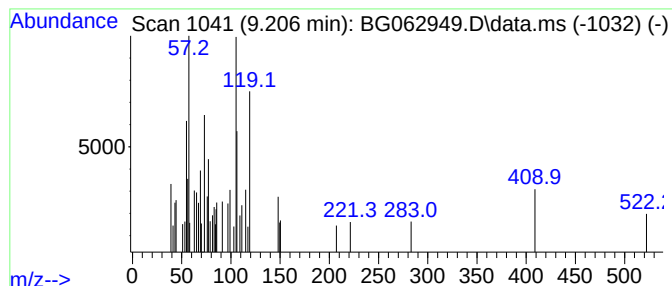
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.206	2.30 ng/ul	31982	1,4-Dichlorobenzene-d4	7.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			piperazin-1-yl(m-tolyl)methanone	204	C12H16N2O	1000455-18-3	25
2			piperazin-1-yl(p-tolyl)methanone	204	C12H16N2O	1000455-18-2	25
3			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	14
4			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	10
5			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
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Misc :
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Instrument :
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ClientSampleId :
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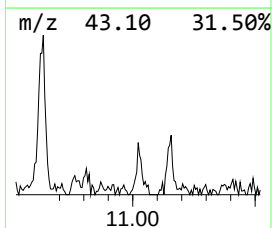
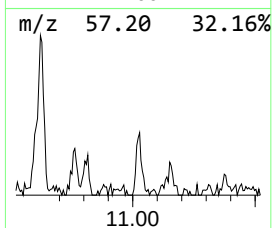
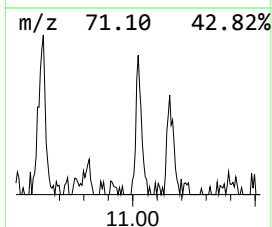
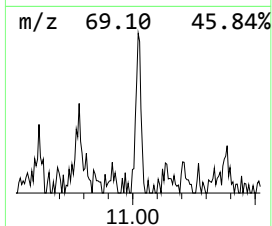
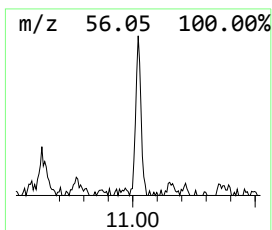
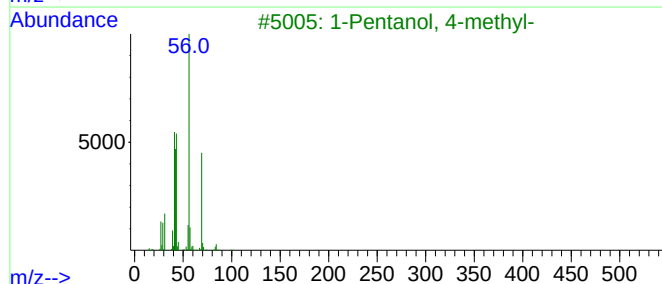
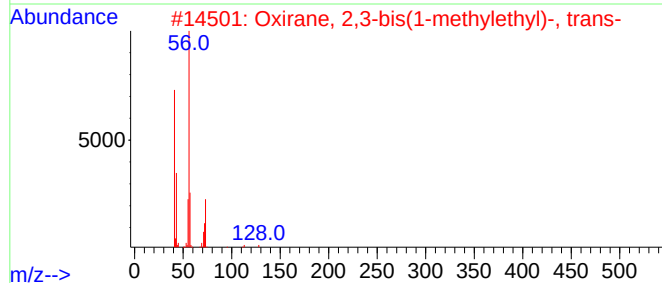
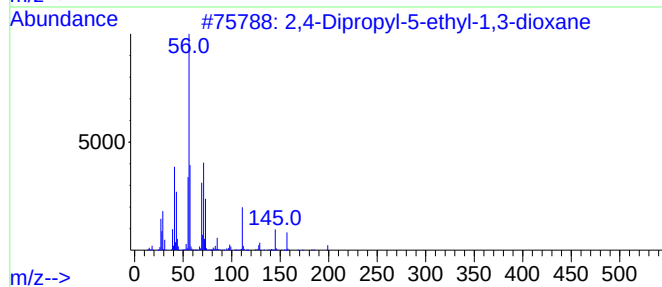
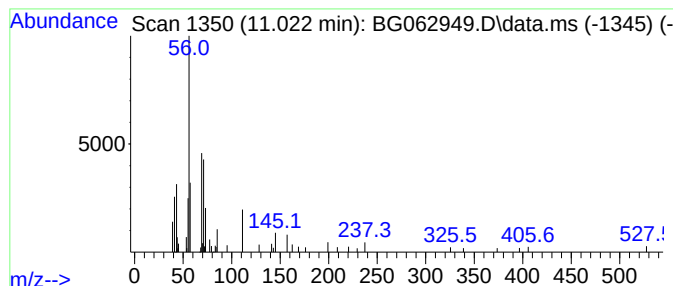
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.022	2.81 ng/ul	59863	Naphthalene-d8	10.640

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	78
2		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	43
3		1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
4		Cyclopentanamine	85	C5H11N	001003-03-8	25
5		Cyclopentanamine	85	C5H11N	001003-03-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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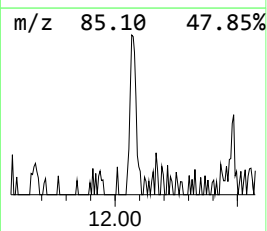
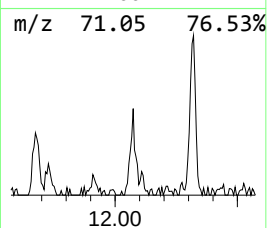
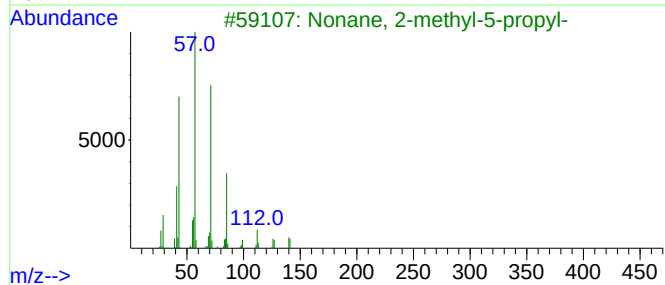
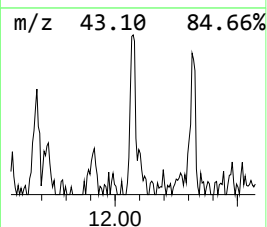
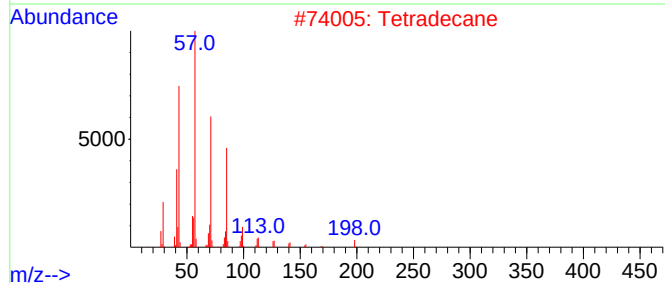
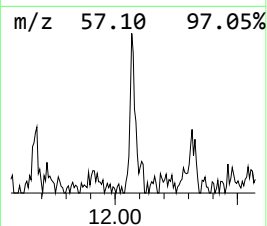
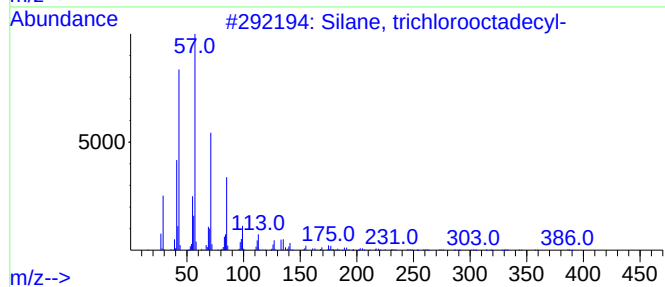
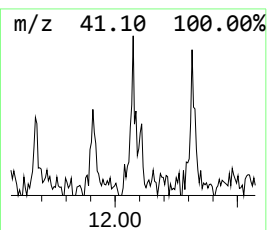
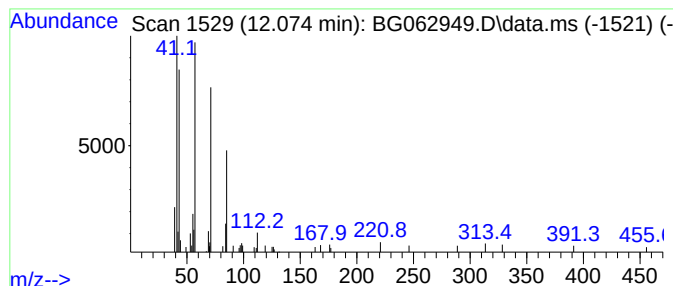
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Silane, trichlorooctadecyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.074	2.29 ng/ul	48834	Naphthalene-d8	10.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	64
2			Tetradecane	198	C14H30	000629-59-4	64
3			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	64
4			Nonane, 5-(1-methylpropyl)-	184	C13H28	062185-54-0	64
5			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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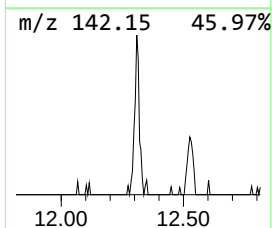
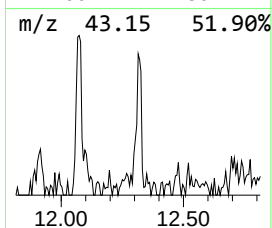
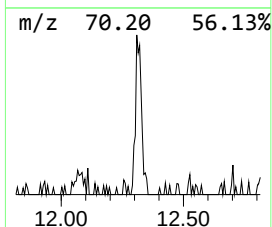
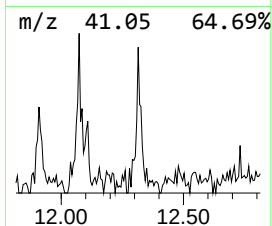
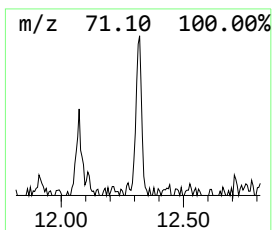
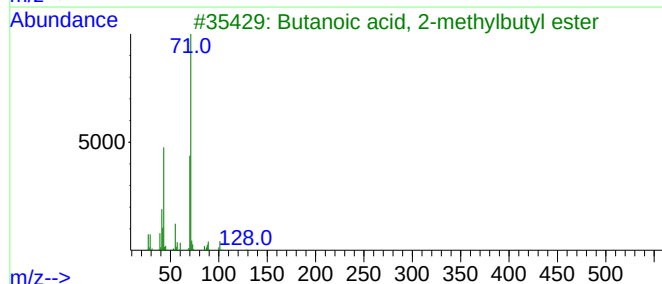
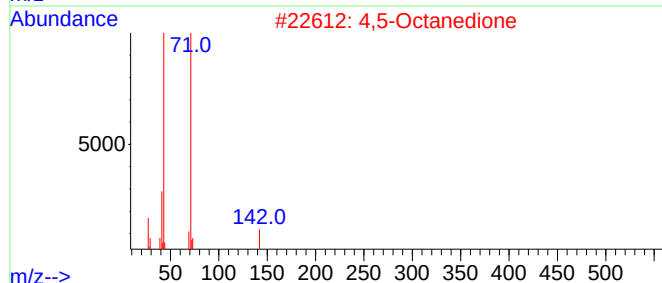
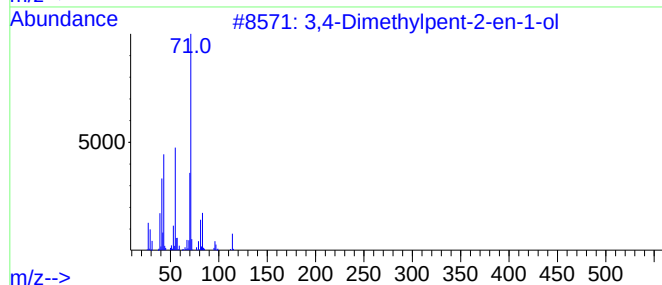
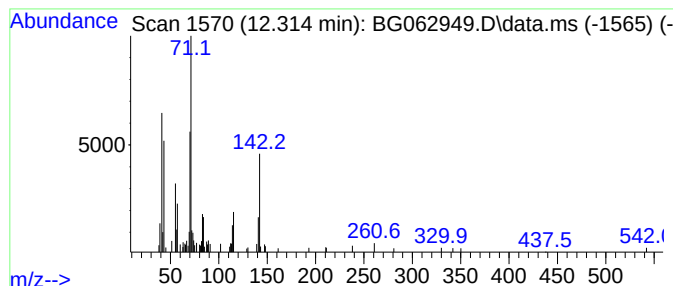
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.315	2.92 ng/ul	62261	Naphthalene-d8	10.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dimethylpent-2-en-1-ol	114	C7H14O	1000195-06-9	38
2			4,5-Octanedione	142	C8H14O2	005455-24-3	38
3			Butanoic acid, 2-methylbutyl ester	158	C9H18O2	051115-64-1	38
4			Propan-1,3-dioic acid, di[3-meth...	244	C13H24O4	064617-96-5	38
5			Dodecane, 4-methyl-	184	C13H28	006117-97-1	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC04MEDL

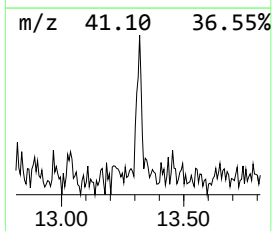
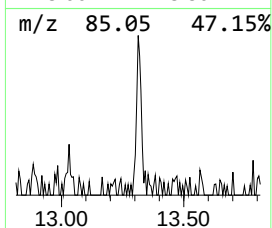
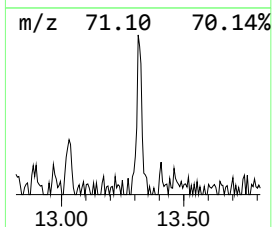
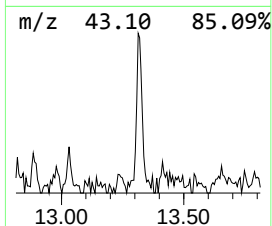
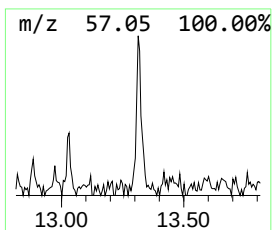
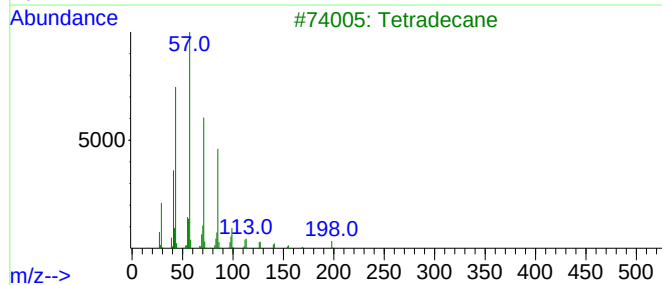
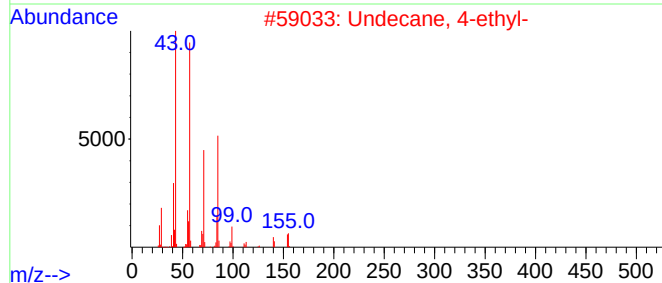
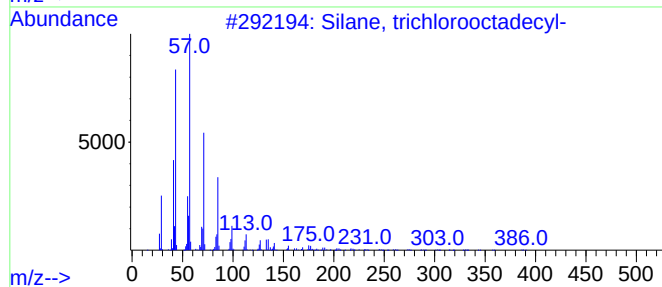
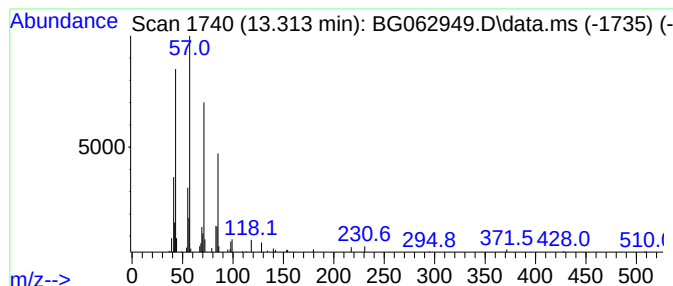
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.313	2.30 ng/ul	59507	Acenaphthene-d10	14.488

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	72
2			Undecane, 4-ethyl-	184	C13H28	017312-59-3	64
3			Tetradecane	198	C14H30	000629-59-4	64
4			Tridecane	184	C13H28	000629-50-5	64
5			Nonane	128	C9H20	000111-84-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

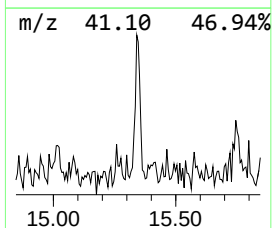
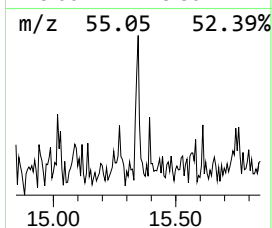
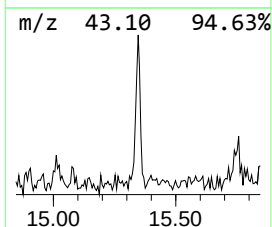
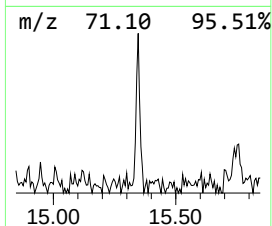
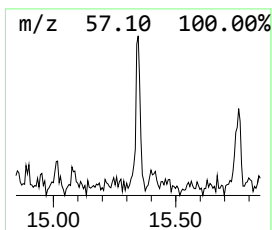
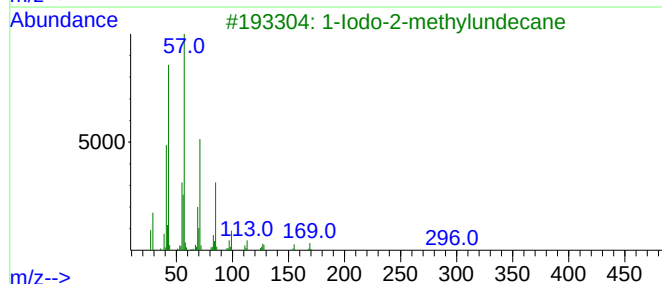
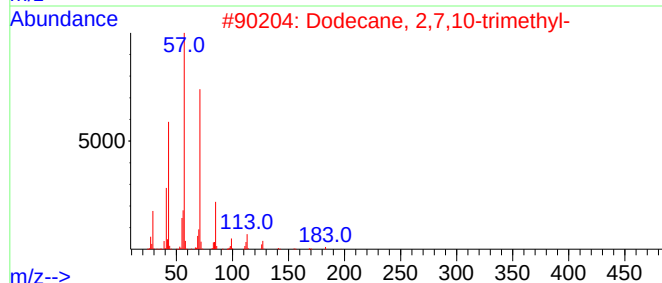
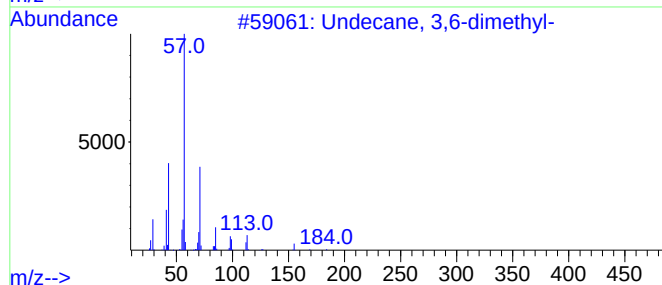
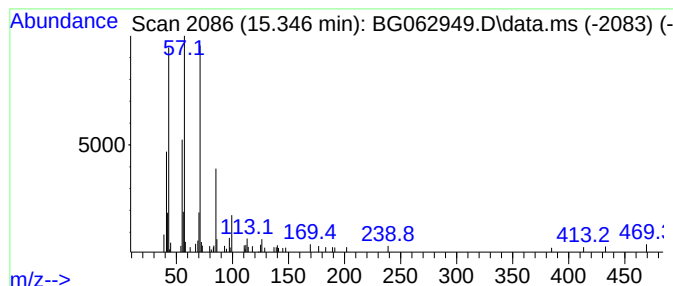
Instrument :
BNA_G
ClientSampleId :
DDC04MEDL

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.346	2.00 ng/ul	51835	Acenaphthene-d10		14.488	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 3,6-dimethyl-		184	C13H28	017301-28-9	72
2	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	59
3	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	53
4	Decane, 2,3,7-trimethyl-		184	C13H28	062238-13-5	50
5	Decane, 3,6-dimethyl-		170	C12H26	017312-53-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062949.D
Acq On : 19 Sep 2024 18:49
Operator : JU/RC
Sample : P3878-15MEDL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC04MEDL

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...	5.887	2.7	ng/u1	36966	1	7.861	277882	20.0
unknown-01	6.856	2.4	ng/u1	33584	1	7.861	277882	20.0
(DEL) Alkane: S...	7.485	11.6	ng/u1	161385	1	7.861	277882	20.0
(DEL) Alkane: S...	7.920	2.7	ng/u1	37974	1	7.861	277882	20.0
1,3-Cyclohexane...	8.284	2.0	ng/u1	27827	1	7.861	277882	20.0
(DEL) Alkane: S...	9.083	8.3	ng/u1	114848	1	7.861	277882	20.0
unknown-02	9.206	2.3	ng/u1	31982	1	7.861	277882	20.0
2,4-Dipropyl-5-...	11.022	2.8	ng/u1	59863	2	10.640	426160	20.0
Silane, trichlo...	12.074	2.3	ng/u1	48834	2	10.640	426160	20.0
unknown-03	12.315	2.9	ng/u1	62261	2	10.640	426160	20.0
(DEL) Alkane: S...	13.313	2.3	ng/u1	59507	3	14.488	518102	20.0
(DEL) Alkane: S...	15.346	2.0	ng/u1	51835	3	14.488	518102	20.0