

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015890.D
 Acq On : 01 Jul 2023 18:47
 Operator : MA/JU
 Sample : 03242-12
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGB3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015890.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	29	32	39	rVB	62807	88269	1.75%	0.114%
2	3.822	105	108	122	rBV	545803	916585	18.22%	1.182%
3	3.922	122	125	134	rVB	66555	99113	1.97%	0.128%
4	4.040	140	145	151	rVB	60765	98986	1.97%	0.128%
5	4.146	160	163	170	rVB	17578	27051	0.54%	0.035%
6	4.722	259	261	265	rVB5	7416	8597	0.17%	0.011%
7	5.093	320	324	326	rBV2	14931	19797	0.39%	0.026%
8	5.122	326	329	336	rVB3	19680	32490	0.65%	0.042%
9	5.228	343	347	348	rBV3	4674	5989	0.12%	0.008%
10	6.693	590	596	605	rVB	64852	107651	2.14%	0.139%
11	7.122	666	669	671	rBV4	4808	5658	0.11%	0.007%
12	7.234	679	688	706	rBV	713050	1800193	35.79%	2.322%
13	7.375	706	712	724	rBV	797757	1371691	27.27%	1.769%
14	7.581	740	747	767	rBV	976111	1818303	36.15%	2.345%
15	8.052	819	827	841	rBV	942102	1798771	35.76%	2.320%
16	8.258	859	862	866	rVB6	5507	8528	0.17%	0.011%
17	8.787	946	952	979	rBV	785050	1979367	39.35%	2.553%
18	9.240	1020	1029	1048	rBV	864420	1817887	36.14%	2.345%
19	9.581	1082	1087	1093	rVB7	8725	18362	0.37%	0.024%
20	9.969	1144	1153	1175	rBV	631204	1548252	30.78%	1.997%
21	10.181	1187	1189	1197	rVB9	9671	19984	0.40%	0.026%
22	10.534	1241	1249	1283	rBV	659488	2221287	44.16%	2.865%
23	10.875	1299	1307	1325	rBV	1366963	2667375	53.03%	3.441%
24	11.051	1327	1337	1362	rBV	454159	1531509	30.45%	1.976%
25	12.040	1498	1505	1508	rBV9	2852	5203	0.10%	0.007%
26	12.081	1508	1512	1518	rVB8	8186	14802	0.29%	0.019%
27	12.428	1562	1571	1579	rBV6	16888	54365	1.08%	0.070%
28	12.498	1579	1583	1591	rVB3	13589	30283	0.60%	0.039%
29	13.075	1677	1681	1686	rVV8	5950	12526	0.25%	0.016%
30	13.304	1715	1720	1722	rBV6	5411	7697	0.15%	0.010%
31	13.498	1749	1753	1757	rBV6	5601	8300	0.17%	0.011%
32	13.575	1760	1766	1776	rVB	25737	49203	0.98%	0.063%
33	13.675	1779	1783	1788	rVB7	9229	15740	0.31%	0.020%
34	14.004	1835	1839	1844	rBV8	6928	12006	0.24%	0.015%
35	14.057	1844	1848	1851	rBV6	3962	6145	0.12%	0.008%
36	14.110	1851	1857	1870	rBV	1914639	3271131	65.03%	4.220%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	14.298	1886	1889	1897	rVB10	4522	5957	0.12%	0.008%
38	14.398	1897	1906	1921	rBV	2343742	3944378	78.42%	5.088%
39	14.563	1930	1934	1940	rVB3	12470	21782	0.43%	0.028%
40	14.698	1950	1957	1978	rVB	2068799	3304991	65.70%	4.263%
41	14.998	2001	2008	2025	rBV2	208156	873032	17.36%	1.126%
42	15.116	2025	2028	2035	rVB6	34218	68366	1.36%	0.088%
43	15.457	2084	2086	2091	rVB6	6868	9186	0.18%	0.012%
44	15.522	2093	2097	2102	rVB6	14813	23084	0.46%	0.030%
45	15.698	2120	2127	2150	rBV	2595470	4632052	92.09%	5.975%
46	15.875	2151	2157	2180	rVV2	305698	878659	17.47%	1.133%
47	16.069	2184	2190	2200	rBV	592184	930585	18.50%	1.200%
48	16.381	2240	2243	2246	rBV4	14623	20361	0.40%	0.026%
49	16.463	2253	2257	2268	rVB3	19705	53772	1.07%	0.069%
50	16.551	2268	2272	2278	rBV9	16283	35093	0.70%	0.045%
51	16.628	2281	2285	2290	rVB	36840	50597	1.01%	0.065%
52	16.845	2318	2322	2326	rBV7	21742	40073	0.80%	0.052%
53	17.175	2375	2378	2382	rBV6	12805	21889	0.44%	0.028%
54	17.339	2402	2406	2412	rBV5	31424	59444	1.18%	0.077%
55	17.463	2421	2427	2438	rBV	2135374	3364067	66.88%	4.339%
56	17.563	2438	2444	2467	rVV2	2271984	4257389	84.64%	5.492%
57	17.733	2470	2473	2486	rVV8	57443	144027	2.86%	0.186%
58	17.839	2488	2491	2496	rVB4	33836	50079	1.00%	0.065%
59	18.204	2549	2553	2555	rBV4	29174	42428	0.84%	0.055%
60	18.245	2555	2560	2566	rBV2	743809	962411	19.13%	1.241%
61	18.316	2567	2572	2580	rBV	1034435	1524778	30.31%	1.967%
62	19.404	2753	2757	2759	rBV	101347	135868	2.70%	0.175%
63	19.539	2776	2780	2785	rBV	238055	364593	7.25%	0.470%
64	19.792	2818	2823	2837	rVB	2897445	5030058	100.00%	6.488%
65	20.263	2900	2903	2913	rVB3	164771	219157	4.36%	0.283%
66	20.751	2983	2986	2988	rBV2	188509	279126	5.55%	0.360%
67	21.222	3060	3066	3076	rVB3	741514	1201689	23.89%	1.550%
68	21.551	3118	3122	3129	rBV2	1624598	2715008	53.98%	3.502%
69	22.121	3217	3219	3224	rVB	320887	382001	7.59%	0.493%
70	23.227	3402	3407	3416	rVB	1087612	1767583	35.14%	2.280%
71	23.927	3521	3526	3539	rVB2	1714105	3869891	76.94%	4.992%
72	24.098	3549	3555	3572	rVB	1423296	3787659	75.30%	4.886%
73	24.551	3625	3632	3641	rVB2	2146516	4682747	93.10%	6.040%
74	24.663	3645	3651	3659	rVB	1127594	2457381	48.85%	3.170%
75	26.615	3978	3983	3993	rVB2	707668	1813395	36.05%	2.339%

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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_P\Methods\SFAM-EPA-BP062623.MA.M
Title : SVOA CALIBRATION

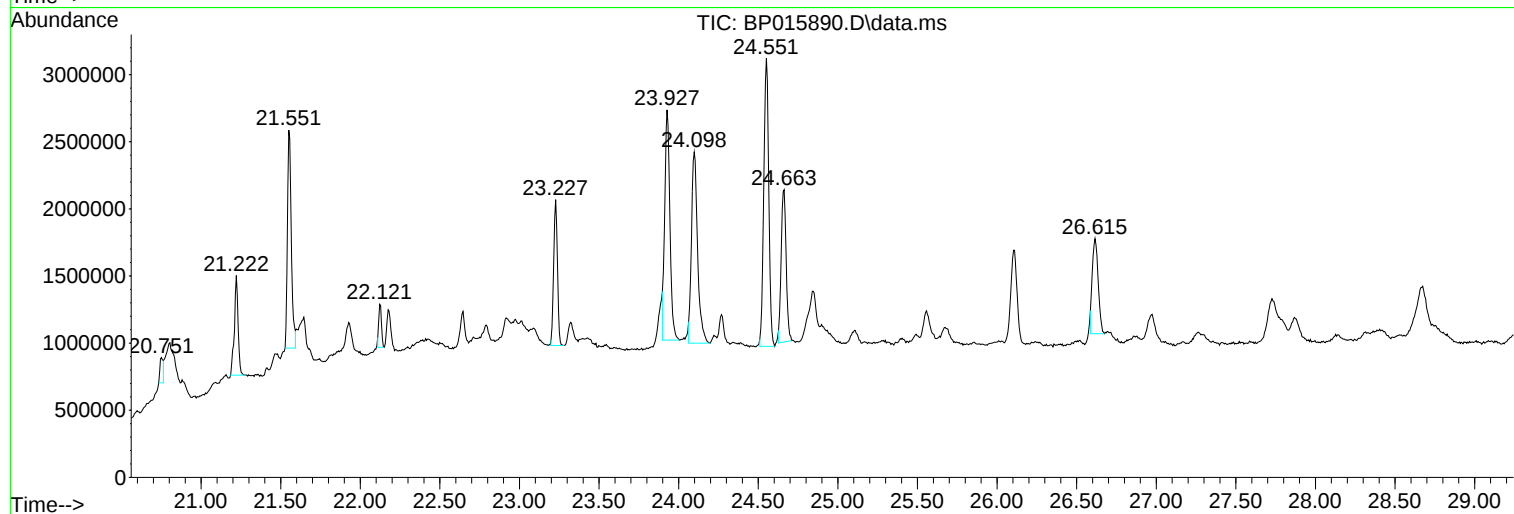
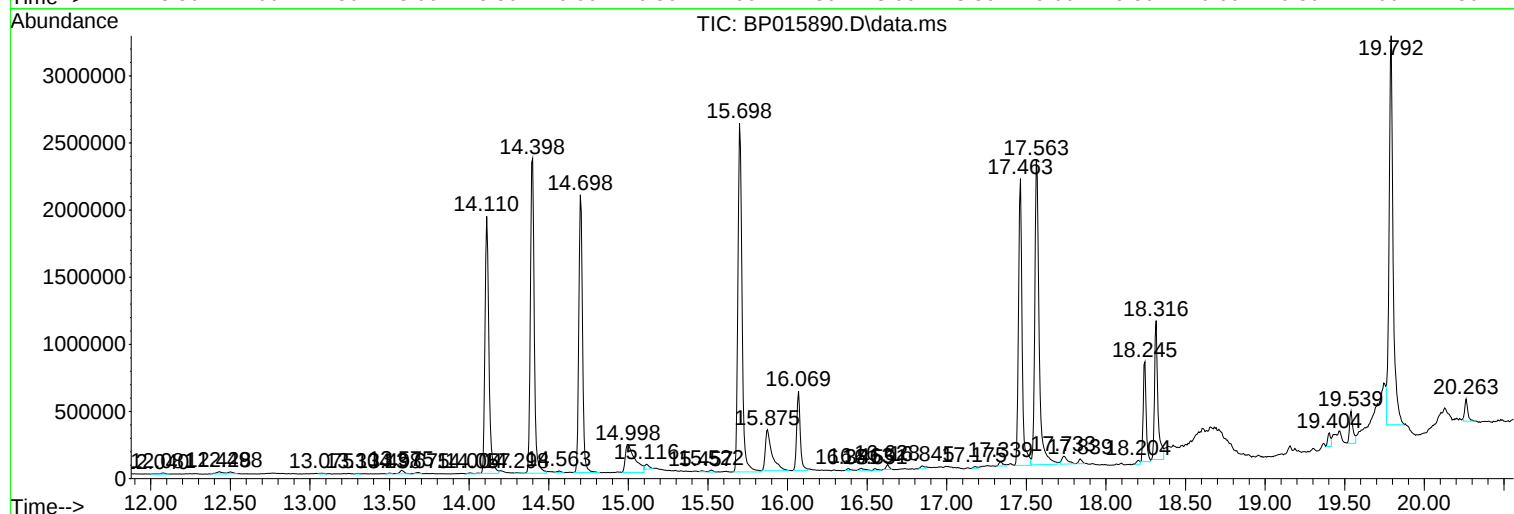
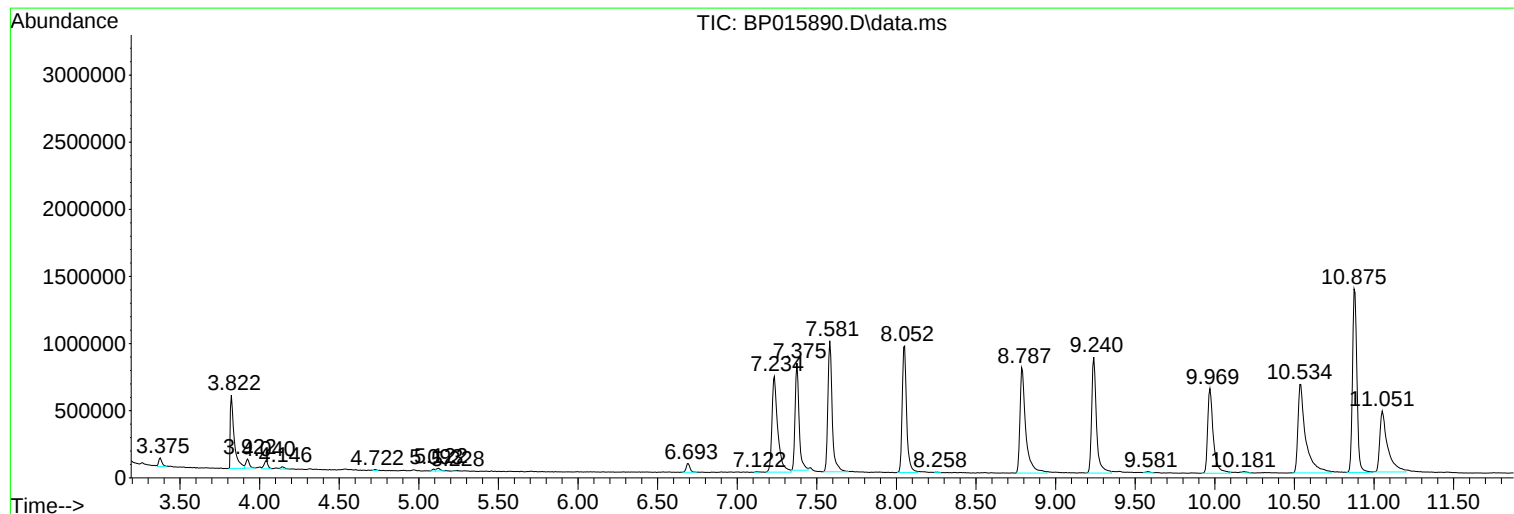
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCGB3

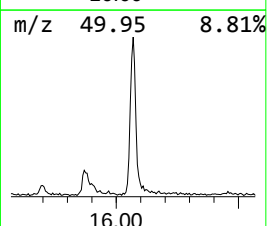
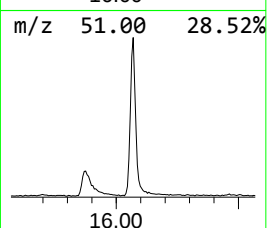
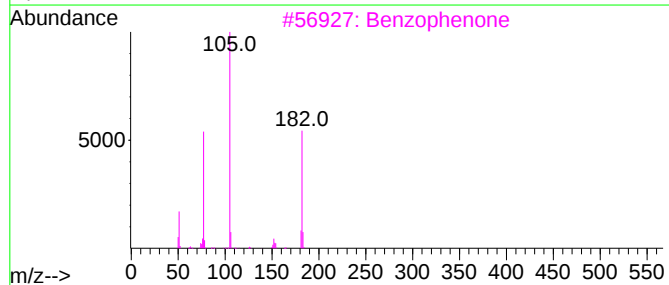
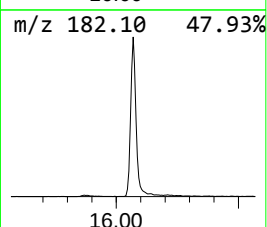
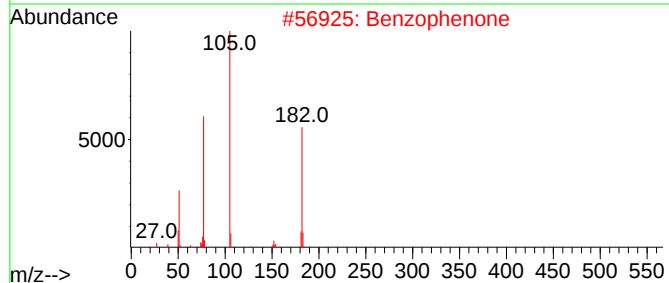
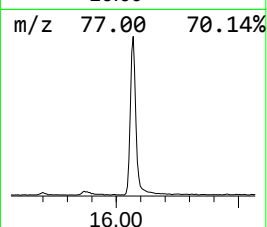
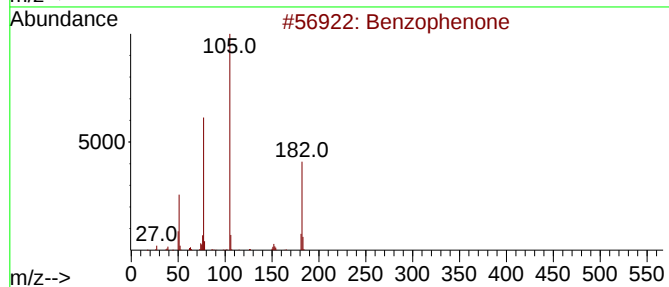
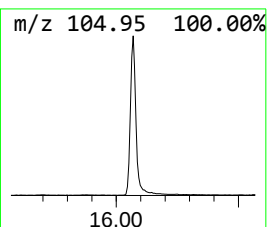
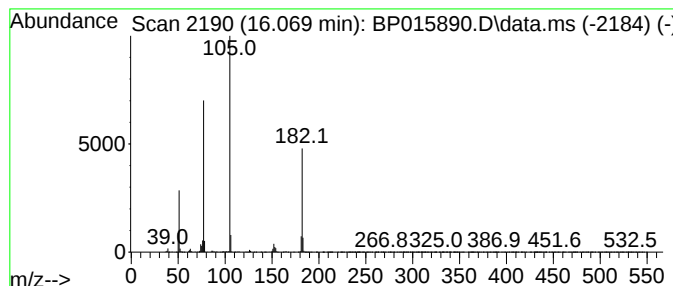
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	5.63 ng/ul	930585	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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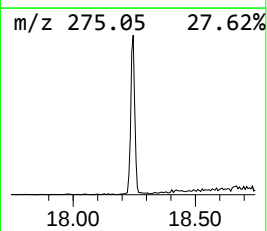
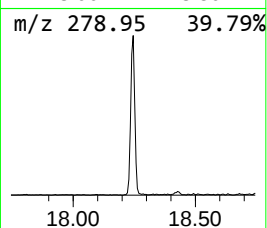
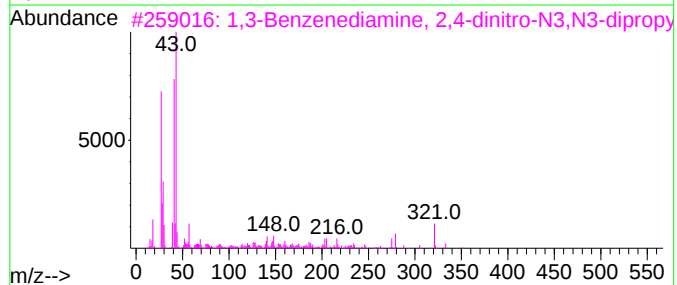
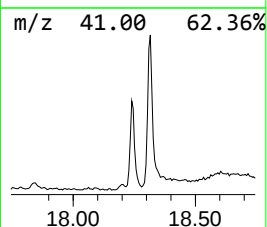
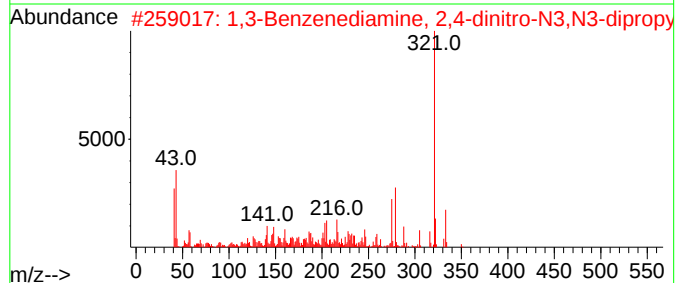
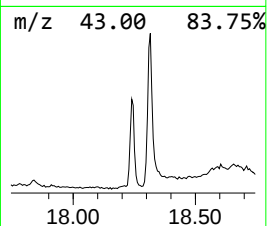
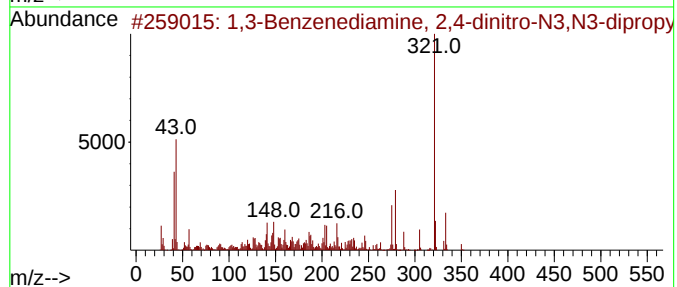
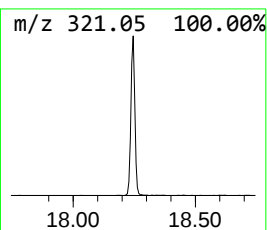
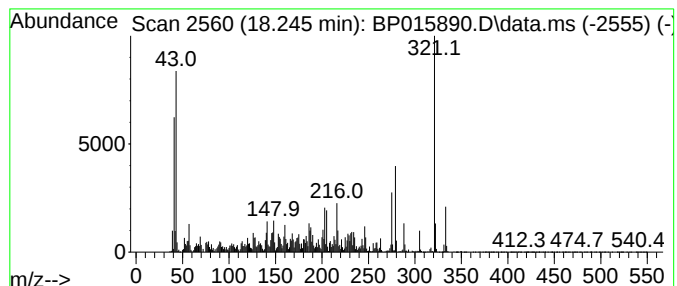
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,3-Benzenediamine, 2,4-din... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	5.72 ng/ul	962411	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	91		
2	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	87		
3	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	68		
4	4-Amino-5-nitro-6-[3,4,5-trimeth...	321	C13H15N5O5	1000254-14-1	38		
5	(E)-4-(2,4,5-Trimethoxystyryl)qu...	321	C20H19NO3	1000442-98-2	16		



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Instrument :
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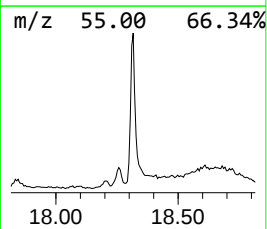
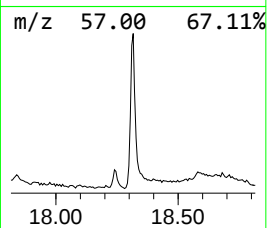
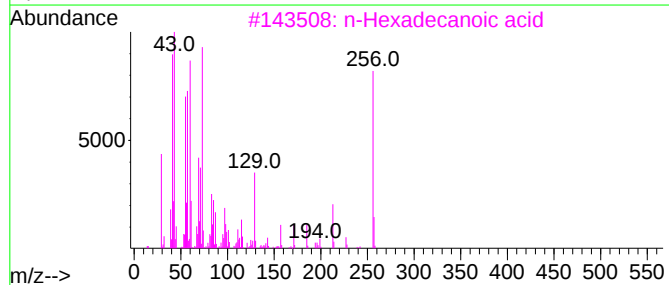
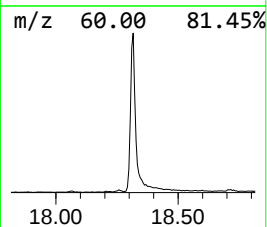
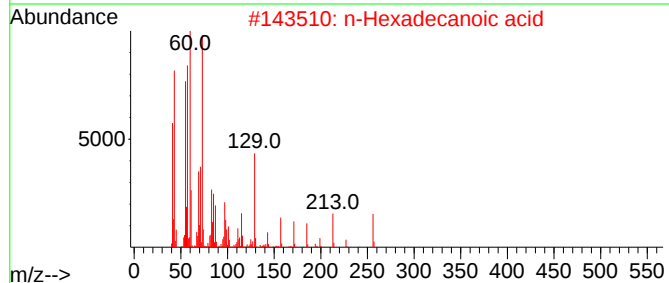
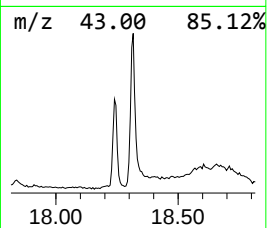
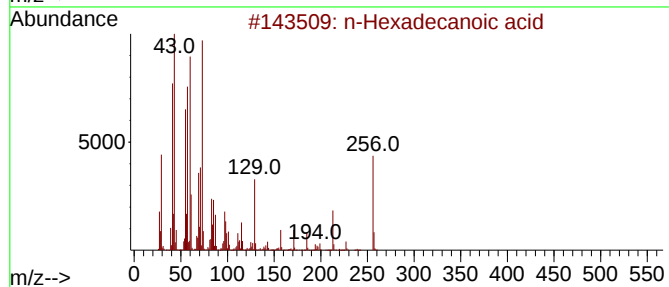
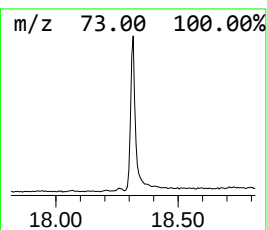
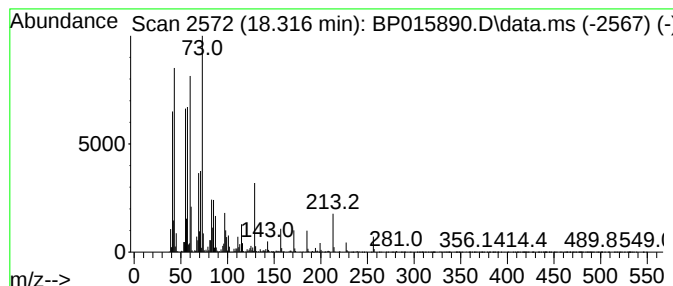
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	9.07 ng/ul	1524780	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97		



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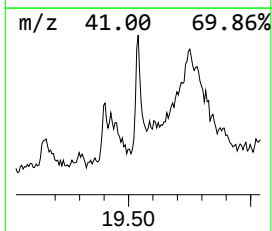
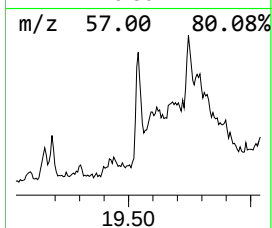
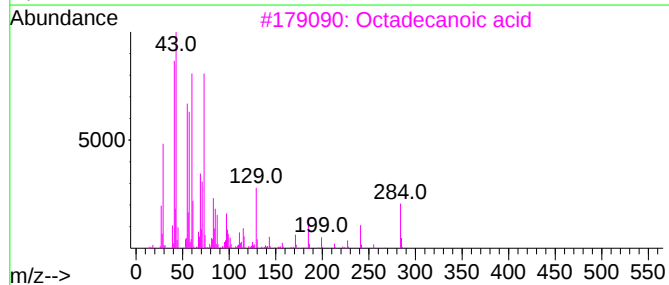
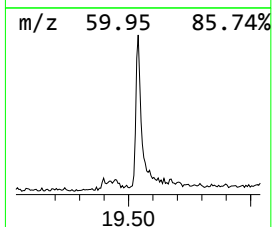
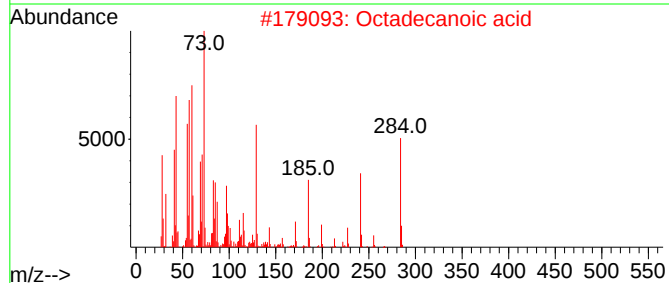
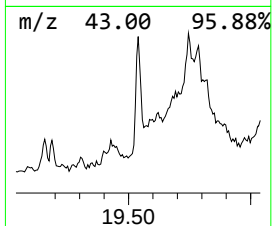
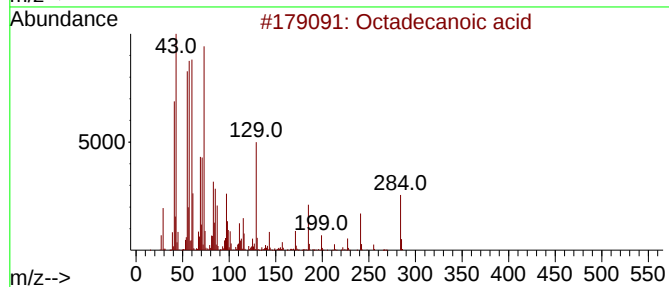
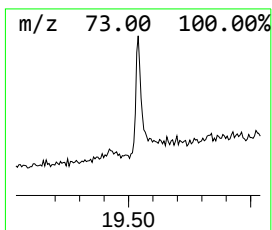
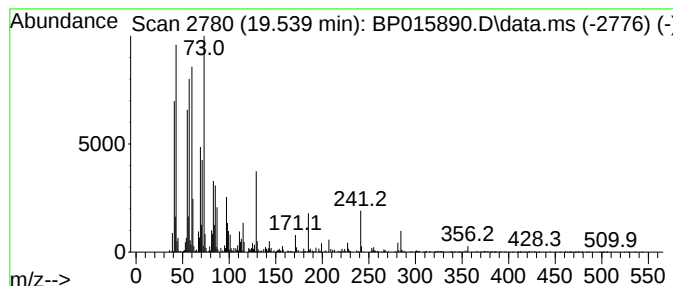
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	2.69 ng/ul	364593	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	96
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
Operator : MA/JU
Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGB3

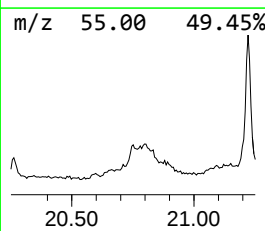
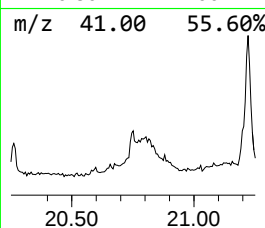
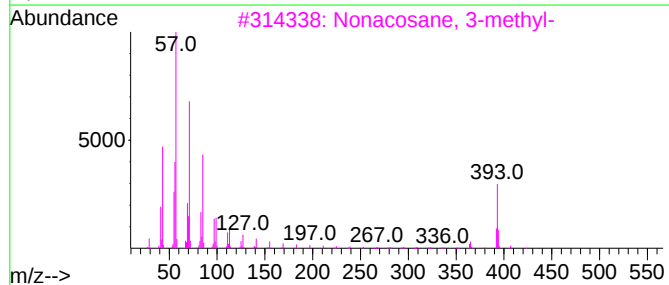
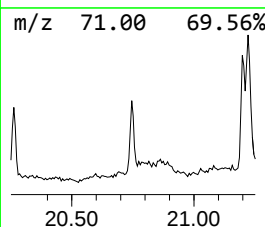
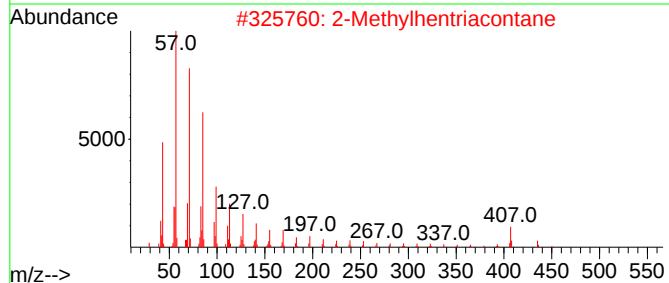
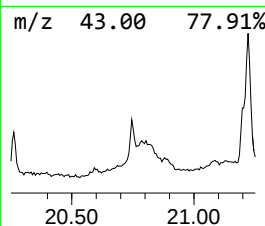
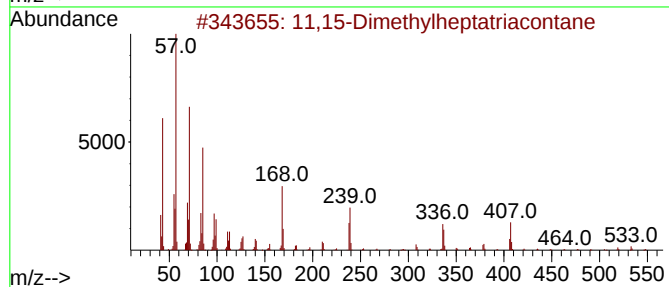
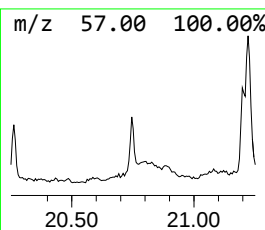
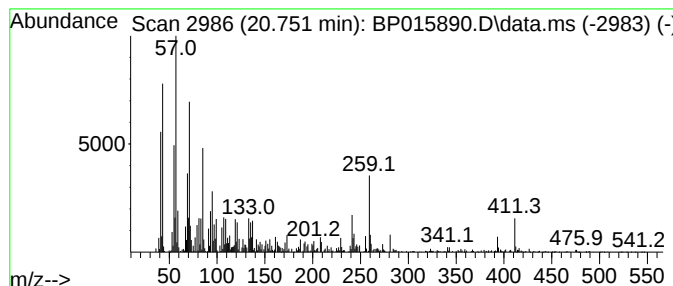
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.751	2.06 ng/ul	279126	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	11,15-Dimethylheptatriacontane	549	C39H80	056987-89-4	32		
2	2-Methylhentriacontane	451	C32H66	001720-12-3	32		
3	Nonacosane, 3-methyl-	422	C30H62	014167-67-0	25		
4	Eicosane, 10-heptyl-10-octyl-	493	C35H72	055470-98-9	22		
5	Isopropyl octacosyl ether	452	C31H64O	1010406-34-7	20		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
Operator : MA/JU
Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

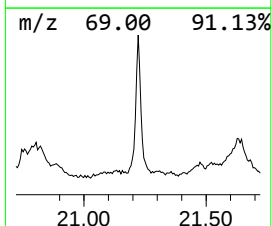
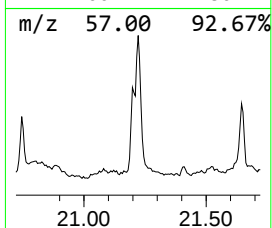
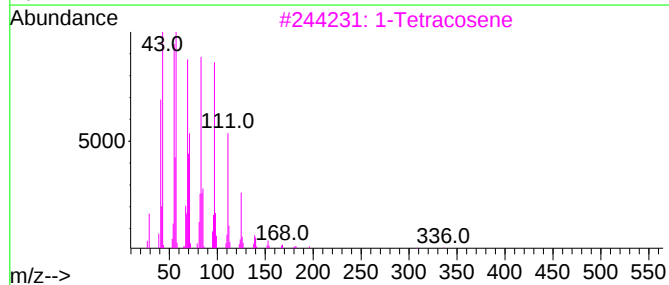
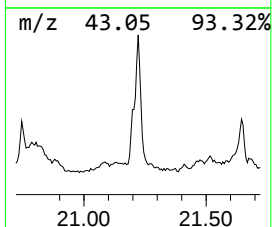
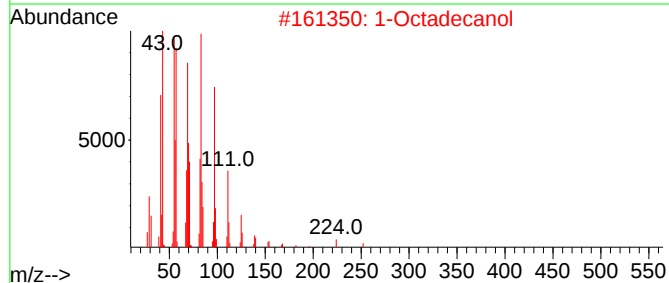
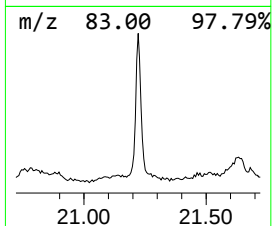
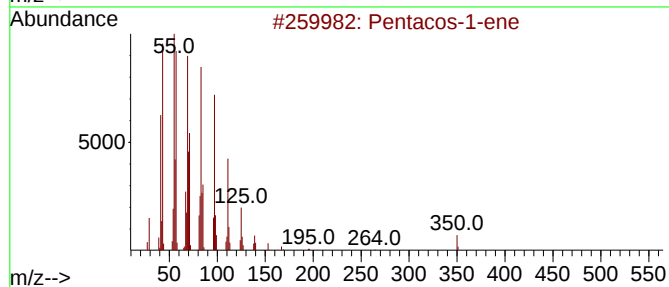
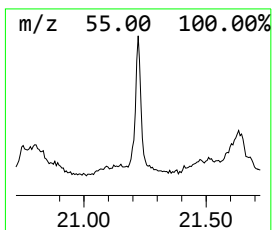
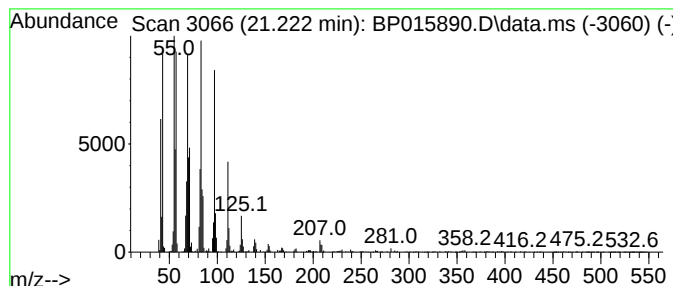
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.221	8.85 ng/ul	1201690	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacos-1-ene	350	C25H50	016980-85-1	94
2	1-Octadecanol	270	C18H38O	000112-92-5	94
3	1-Tetracosene	336	C24H48	010192-32-2	94
4	1-Hexadecanol	242	C16H34O	036653-82-4	93
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
Operator : MA/JU
Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

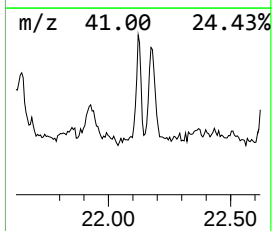
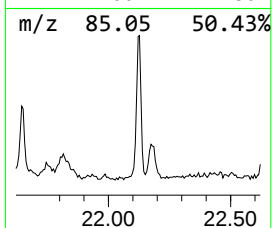
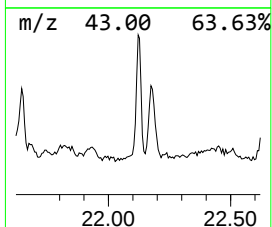
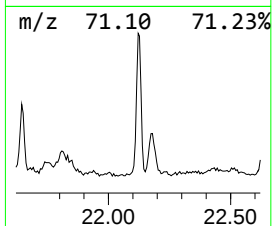
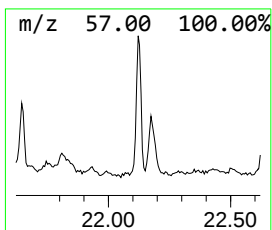
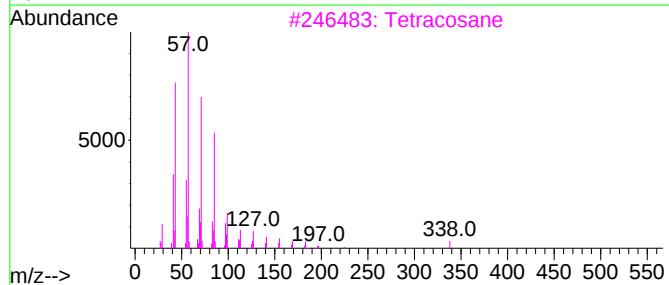
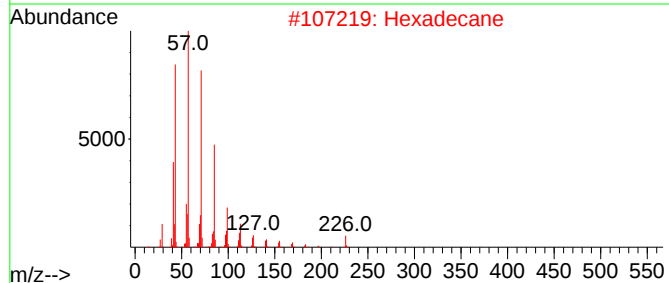
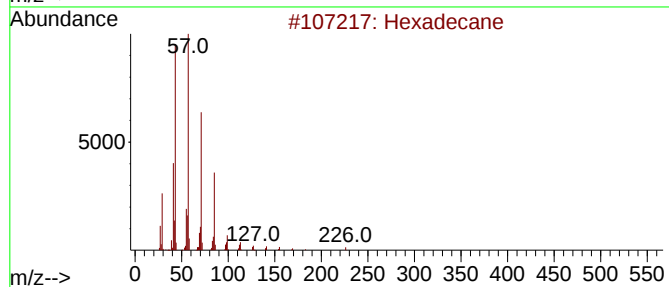
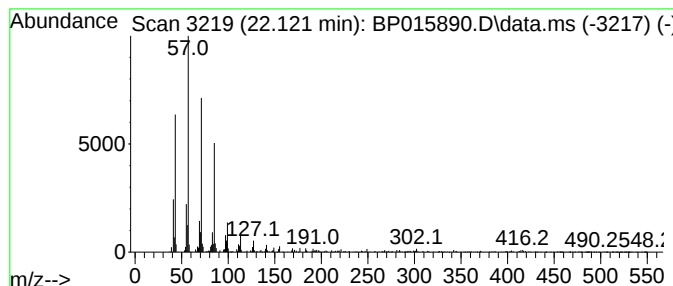
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.122	2.81 ng/ul	382001	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Hexadecane	226	C16H34	000544-76-3	87
3	Tetracosane	338	C24H50	000646-31-1	87
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5	Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
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Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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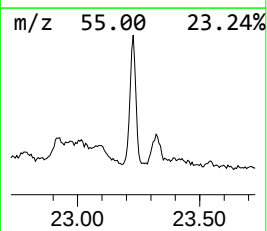
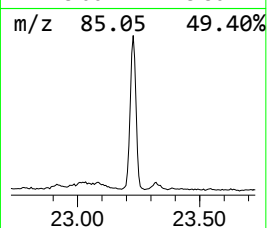
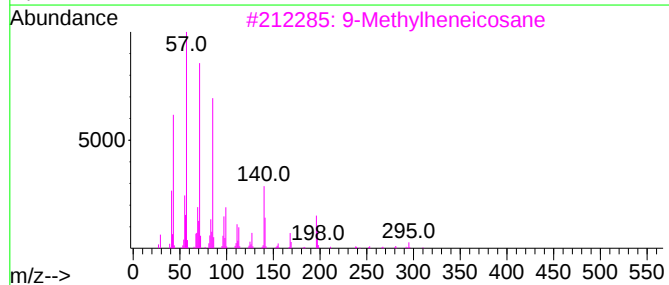
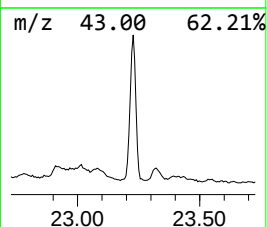
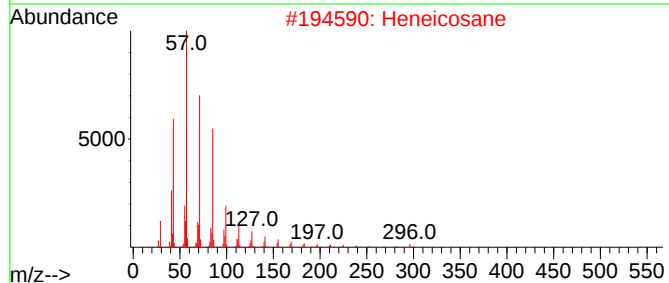
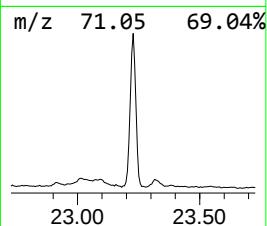
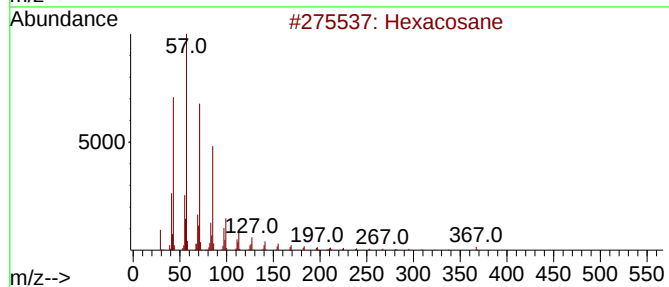
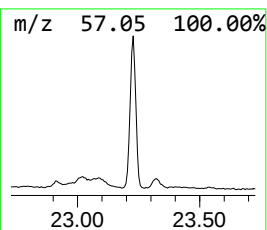
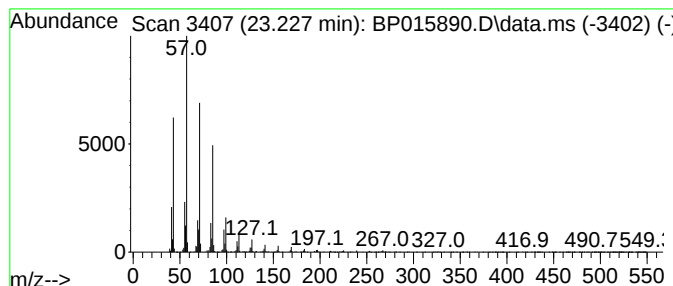
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	9.33 ng/ul	1767580	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			9-Methylheneicosane	310	C22H46	070475-51-3	91
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5			Eicosane	282	C20H42	000112-95-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
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Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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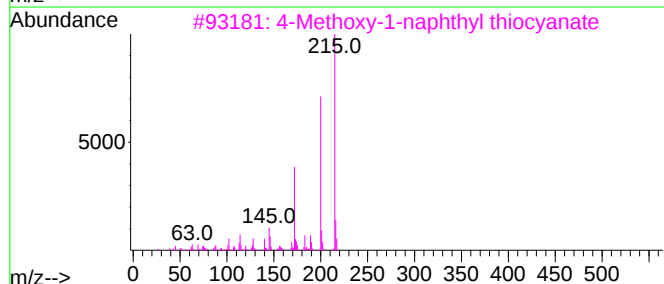
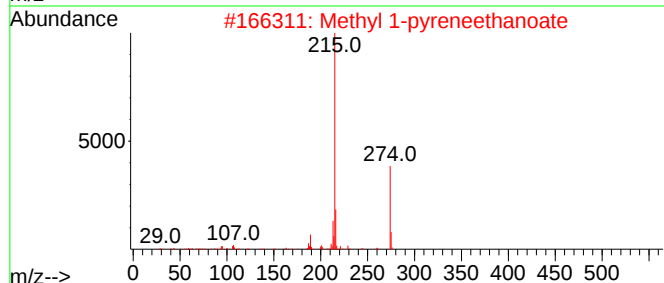
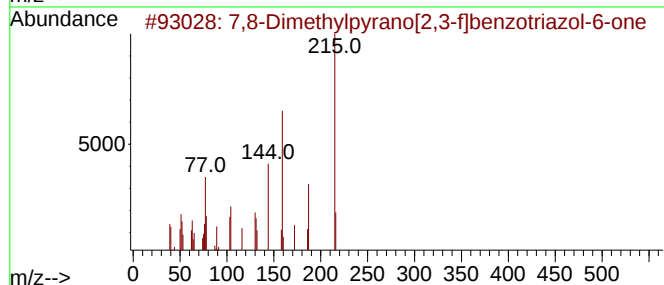
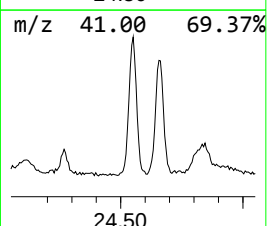
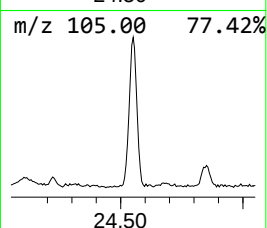
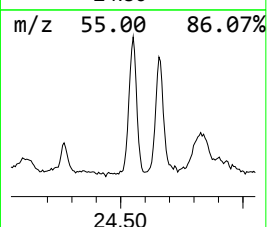
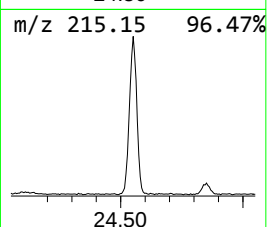
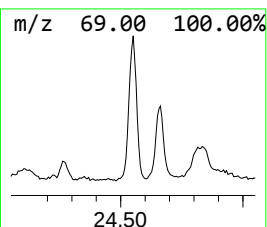
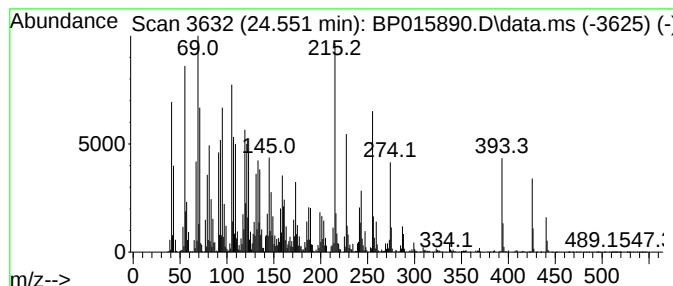
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.551	24.73 ng/ul	4682750	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,8-Dimethylpyrano[2,3-f]benzotr...	215	C11H9N3O2	129503-06-6	25
2			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	20
3			4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	15
4			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	15
5			Indanidine	215	C11H13N5	085392-79-6	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
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Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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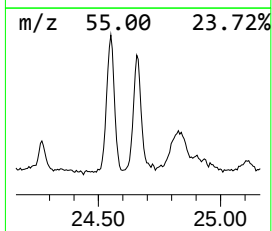
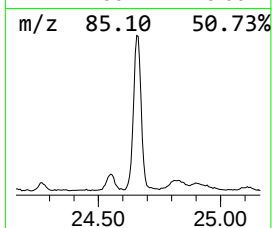
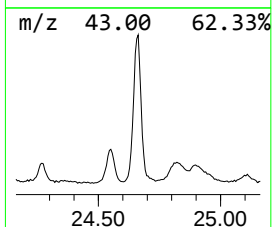
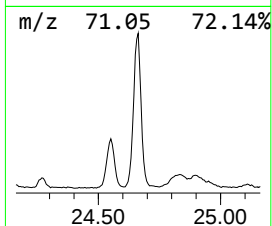
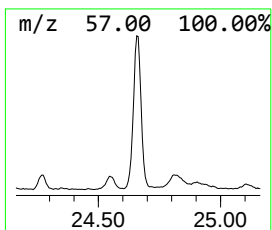
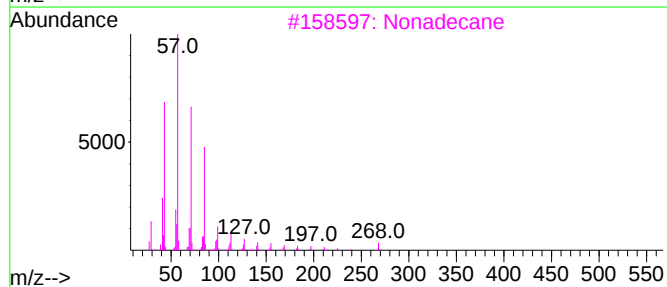
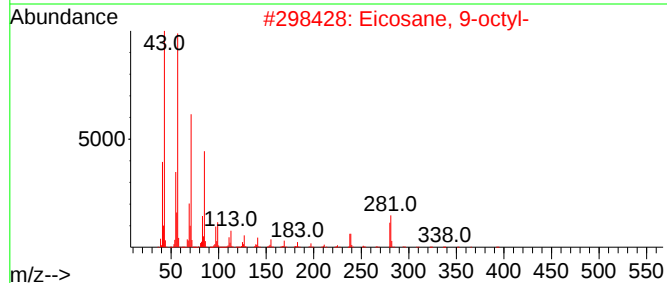
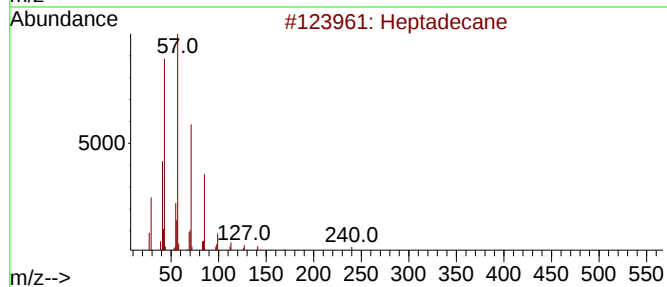
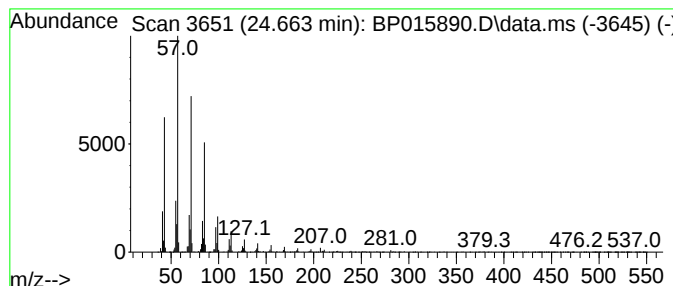
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.663	12.98 ng/ul	2457380	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
3		Nonadecane	268	C19H40	000629-92-5	93
4		2-Methylheptacosane	394	C28H58	001561-00-8	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015890.D
Acq On : 01 Jul 2023 18:47
Operator : MA/JU
Sample : 03242-12
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGB3

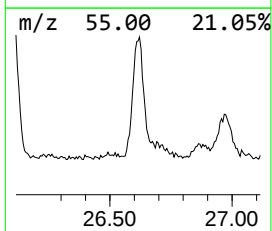
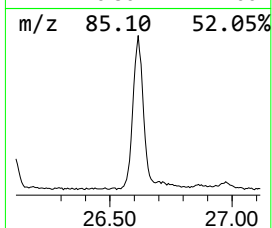
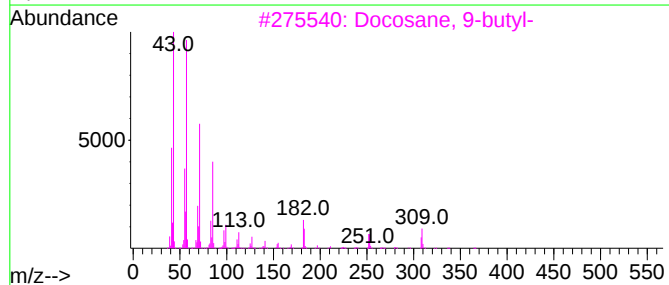
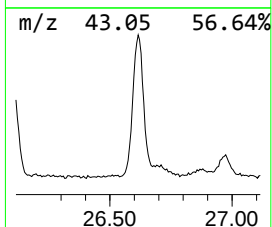
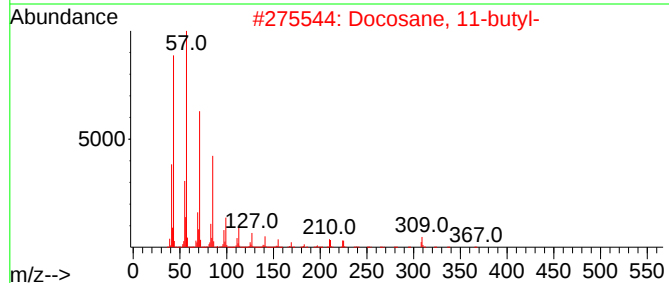
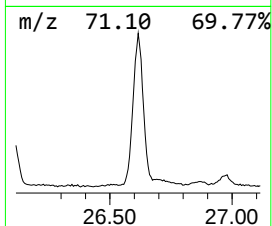
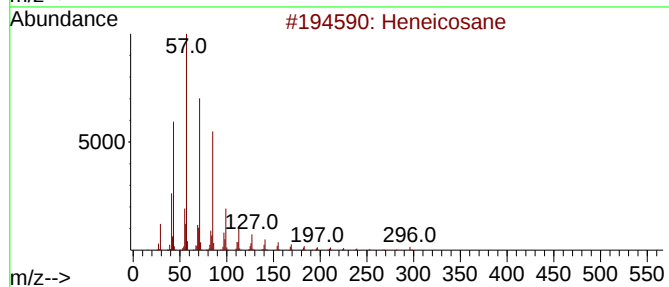
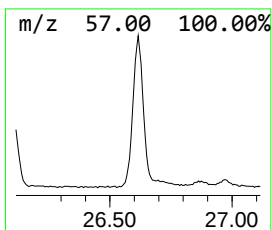
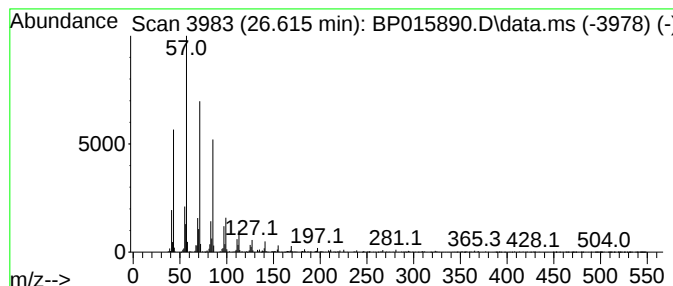
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.615	9.58 ng/ul	1813400	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	98
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	95
3			Docosane, 9-butyl-	366	C26H54	055282-14-9	94
4			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
5			Docosane, 5-butyl-	366	C26H54	055282-16-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015890.D
 Acq On : 01 Jul 2023 18:47
 Operator : MA/JU
 Sample : 03242-12
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGB3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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
Benzophenone	16.069	5.6	ng/ul	930585	3	14.698	3304990	20.0
1,3-Benzenediam...	18.245	5.7	ng/ul	962411	4	17.463	3364070	20.0
n-Hexadecanoic ...	18.316	9.1	ng/ul	1524780	4	17.463	3364070	20.0
Octadecanoic acid	19.539	2.7	ng/ul	364593	5	21.551	2715010	20.0
(DEL) Alkane: S...	20.751	2.1	ng/ul	279126	5	21.551	2715010	20.0
(DEL) Alkane: S...	21.221	8.8	ng/ul	1201690	5	21.551	2715010	20.0
(DEL) Alkane: S...	22.122	2.8	ng/ul	382001	5	21.551	2715010	20.0
(DEL) Alkane: S...	23.227	9.3	ng/ul	1767580	6	24.098	3787660	20.0
unknown-01	24.551	24.7	ng/ul	4682750	6	24.098	3787660	20.0
(DEL) Alkane: S...	24.663	13.0	ng/ul	2457380	6	24.098	3787660	20.0
(DEL) Alkane: S...	26.615	9.6	ng/ul	1813400	6	24.098	3787660	20.0