

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
 Data File : BM024508.D
 Acq On : 17 Jan 2020 22:46
 Operator : JU
 Sample : L1109-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GASG6

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_M\METHODS\SOM-EPA-BM011520MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.063	26	30	35	rVB	45459	58509	0.92%	0.080%
2	3.334	73	76	81	rBV2	9259	15524	0.24%	0.021%
3	3.681	132	135	140	rVB3	8898	12592	0.20%	0.017%
4	3.763	146	149	156	rVB5	7611	12452	0.20%	0.017%
5	4.646	296	299	306	rVB5	7737	12600	0.20%	0.017%
6	6.646	633	639	651	rBV	784139	1294750	20.29%	1.781%
7	6.804	660	666	675	rBV	578651	902517	14.14%	1.242%
8	6.998	692	699	711	rBV	897960	1367795	21.43%	1.882%
9	7.457	771	777	787	rBV	787357	1211994	18.99%	1.667%
10	7.545	787	792	798	rVB	100607	144104	2.26%	0.198%
11	8.169	891	898	909	rBV	1103355	1706399	26.74%	2.347%
12	8.592	963	970	983	rBV	875696	1417810	22.22%	1.950%
13	9.098	1051	1056	1061	rVB	18592	27414	0.43%	0.038%
14	9.310	1085	1092	1103	rBV	792687	1259918	19.74%	1.733%
15	9.392	1103	1106	1112	rVB5	7910	10340	0.16%	0.014%
16	9.845	1175	1183	1194	rBV	1204873	1948070	30.53%	2.680%
17	10.216	1239	1246	1254	rBV	1125654	1819998	28.52%	2.504%
18	10.351	1263	1269	1287	rBV	1101308	1851973	29.02%	2.548%
19	11.816	1513	1518	1523	rVB2	18381	28405	0.45%	0.039%
20	13.345	1773	1778	1786	rBV	91787	127672	2.00%	0.176%
21	13.527	1802	1809	1813	rBV	2447583	3266903	51.20%	4.494%
22	13.574	1813	1817	1823	rVB	558587	727071	11.39%	1.000%
23	13.792	1847	1854	1864	rBV	2647688	3748906	58.75%	5.157%
24	14.104	1901	1907	1918	rVB	1916721	2783060	43.61%	3.829%
25	14.316	1936	1943	1960	rBV	1007489	1517708	23.78%	2.088%
26	14.539	1977	1981	1985	rBV5	18857	26738	0.42%	0.037%
27	14.957	2046	2052	2056	rBV6	6585	9133	0.14%	0.013%
28	15.016	2056	2062	2065	rBV6	7649	9072	0.14%	0.012%
29	15.104	2071	2077	2087	rBV	3476442	4792905	75.11%	6.594%
30	15.227	2092	2098	2112	rVV	973331	1399987	21.94%	1.926%
31	15.345	2114	2118	2124	rVB	37422	52091	0.82%	0.072%
32	15.568	2153	2156	2164	rVB6	6694	9621	0.15%	0.013%
33	15.892	2206	2211	2217	rBV5	12834	22263	0.35%	0.031%
34	16.298	2275	2280	2283	rBV4	7592	10912	0.17%	0.015%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
 Data File : BM024508.D
 Acq On : 17 Jan 2020 22:46
 Operator : JU
 Sample : L1109-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GASG6

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_M\METHODS\SOM-EPA-BM011520MA.M
 Title : SVOA CALIBRATION

35	16.539	2316	2321	2324	rVB4	5413	7116	0.11%	0.010%
36	16.645	2334	2339	2341	rBV	9253	13155	0.21%	0.018%
37	16.674	2341	2344	2347	rVB4	7604	9152	0.14%	0.013%
38	16.851	2365	2374	2385	rBV	2453438	3523847	55.22%	4.848%
39	16.951	2385	2391	2403	rVV	3716842	5250550	82.28%	7.223%
40	17.068	2407	2411	2415	rVB7	15865	20220	0.32%	0.028%
41	17.280	2442	2447	2452	rBV4	10727	15245	0.24%	0.021%
42	17.415	2466	2470	2473	rVB4	5376	7833	0.12%	0.011%
43	17.668	2504	2513	2514	rBV8	5297	10549	0.17%	0.015%
44	17.798	2529	2535	2553	rBV	680837	1020648	15.99%	1.404%
45	18.468	2645	2649	2653	rBV6	8443	12407	0.19%	0.017%
46	18.609	2669	2673	2677	rVB6	8935	14649	0.23%	0.020%
47	18.668	2677	2683	2688	rBV9	8267	19582	0.31%	0.027%
48	18.745	2693	2696	2699	rVB5	9458	10283	0.16%	0.014%
49	18.886	2716	2720	2725	rBV	40477	55440	0.87%	0.076%
50	18.939	2725	2729	2731	rVV4	24773	36992	0.58%	0.051%
51	18.962	2731	2733	2742	rVB6	41406	65368	1.02%	0.090%
52	19.092	2750	2755	2760	rBV	344090	482156	7.56%	0.663%
53	19.139	2760	2763	2772	rVB	55865	88661	1.39%	0.122%
54	19.256	2777	2783	2793	rVV	4873921	6381158	100.00%	8.779%
55	19.327	2793	2795	2798	rVV3	20020	26204	0.41%	0.036%
56	19.362	2798	2801	2805	rVB6	13552	17455	0.27%	0.024%
57	19.827	2877	2880	2884	rBV3	38341	47192	0.74%	0.065%
58	19.868	2884	2887	2893	rVB6	34847	45350	0.71%	0.062%
59	20.180	2938	2940	2946	rBV6	13054	25842	0.40%	0.036%
60	20.227	2946	2948	2952	rVV5	14222	19647	0.31%	0.027%
61	20.368	2969	2972	2975	rVB	59510	59107	0.93%	0.081%
62	20.815	3044	3048	3061	rBV	1190855	1654361	25.93%	2.276%
63	21.056	3084	3089	3098	rBV	3316275	4191163	65.68%	5.766%
64	21.268	3121	3125	3129	rBV	298484	316447	4.96%	0.435%
65	21.492	3160	3163	3166	rBV3	68929	69732	1.09%	0.096%
66	21.692	3193	3197	3204	rBV	505981	602143	9.44%	0.828%
67	22.133	3268	3272	3284	rVB	571581	761222	11.93%	1.047%
68	22.244	3288	3291	3297	rVB	186468	234291	3.67%	0.322%
69	22.609	3349	3353	3361	rVB	642660	870388	13.64%	1.197%
70	23.044	3420	3427	3437	rVB	3689320	6117656	95.87%	8.416%
71	23.144	3439	3444	3445	rVV	646585	882851	13.84%	1.215%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

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_M\METHODS\SOM-EPA-BM011520MA.M
Title : SVOA CALIBRATION

72	23.180	3445	3450	3461	rVB2	2300412	4056424	63.57%	5.580%
73	23.744	3541	3546	3552	rBV	488957	841064	13.18%	1.157%
74	23.809	3553	3557	3569	rVB2	285938	560315	8.78%	0.771%
75	24.438	3659	3664	3672	rVB	322351	639339	10.02%	0.880%

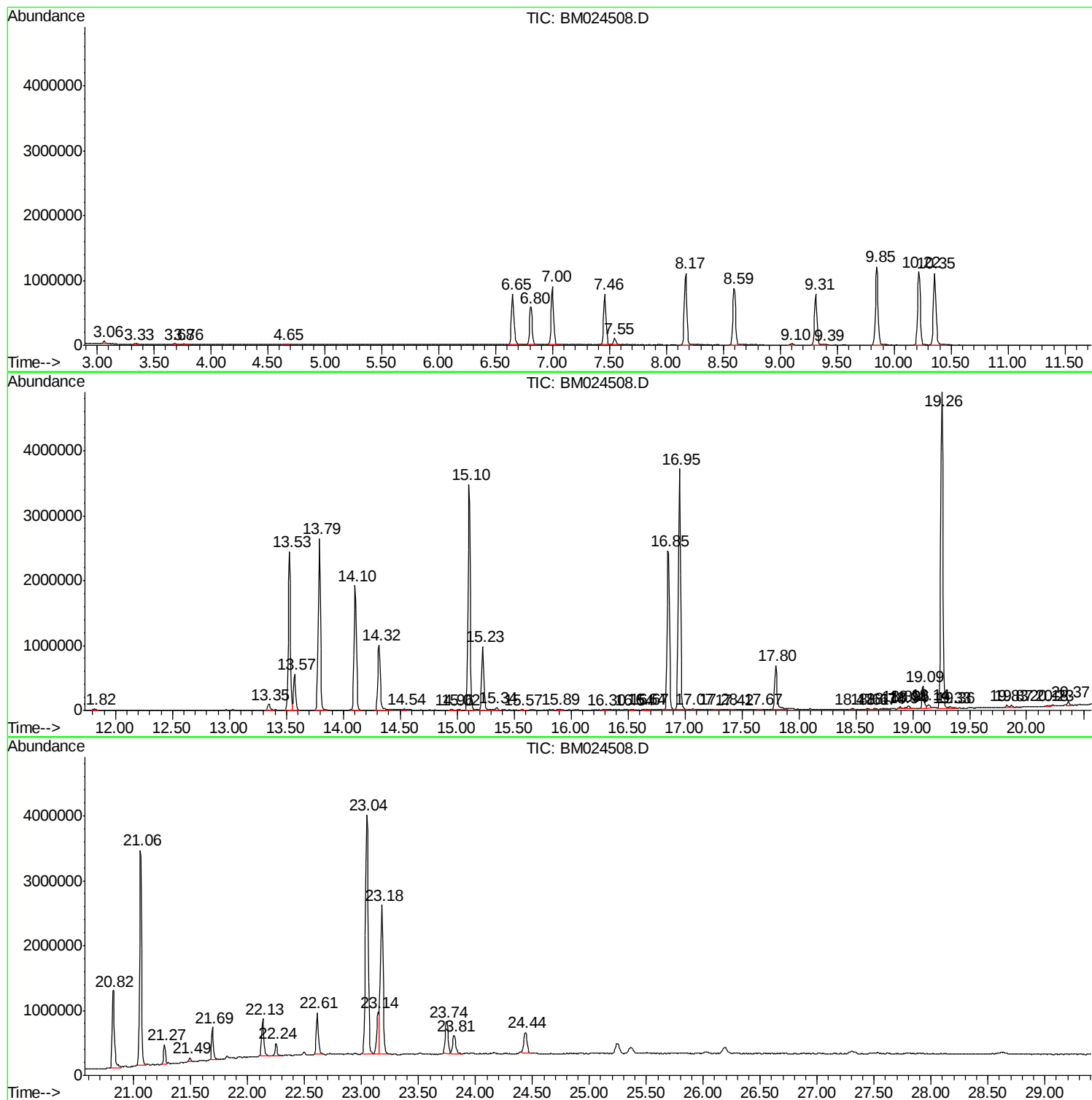
Sum of corrected areas: 72690410

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

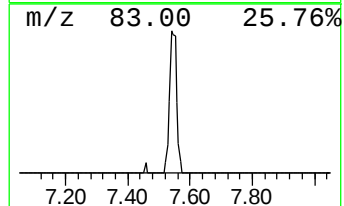
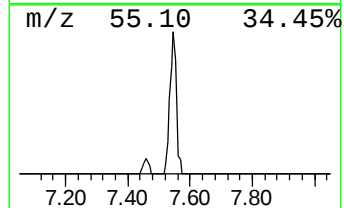
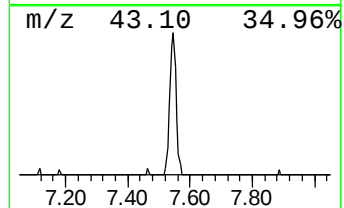
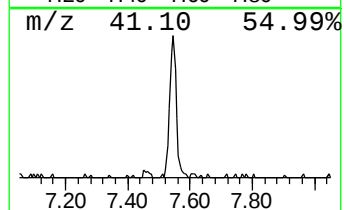
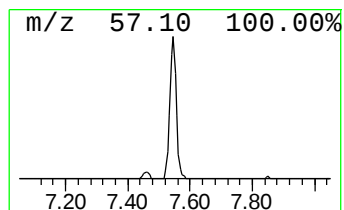
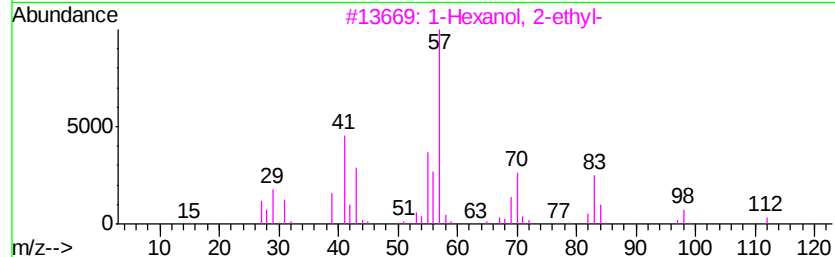
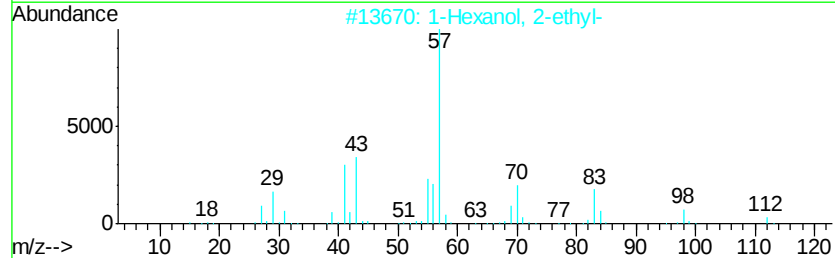
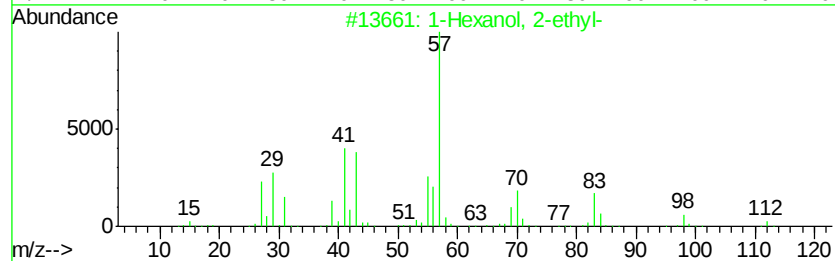
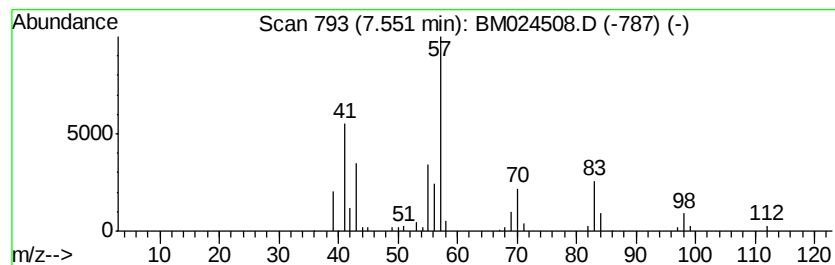
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	2.38 ng/ul	144104	1,4-Dichlorobenzene-d4	7.46

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
5			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

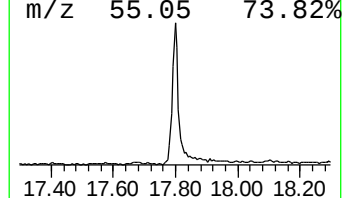
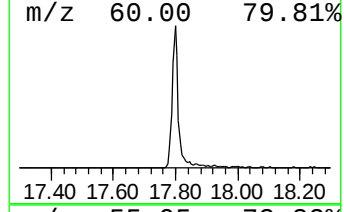
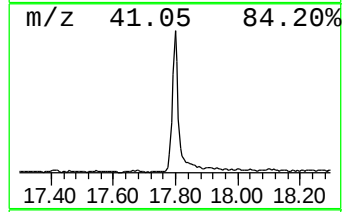
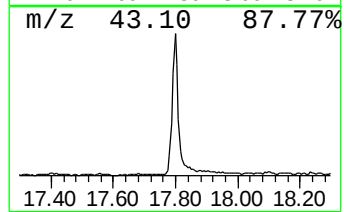
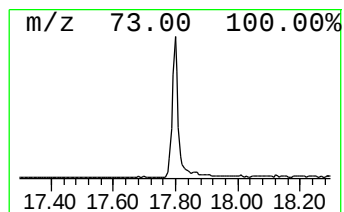
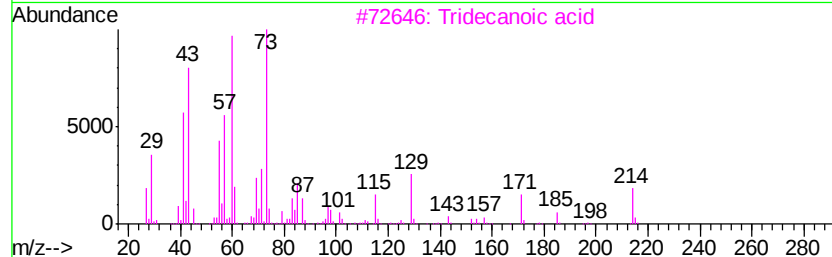
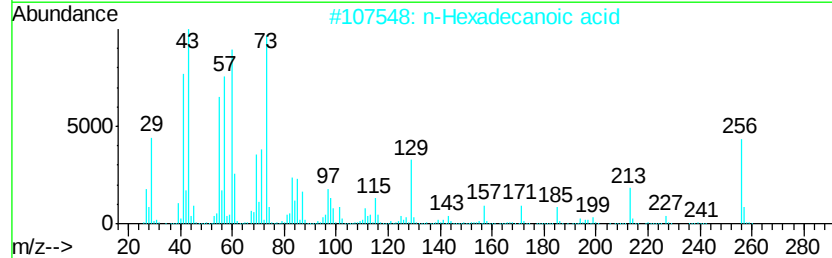
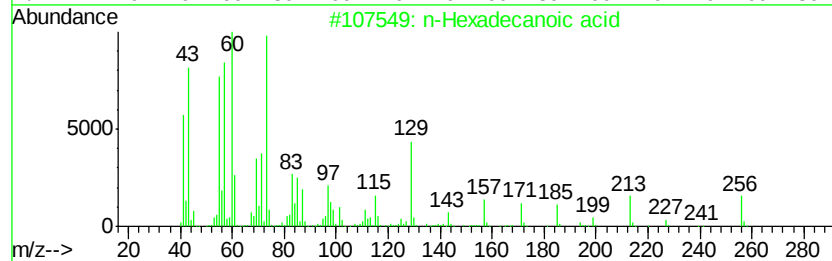
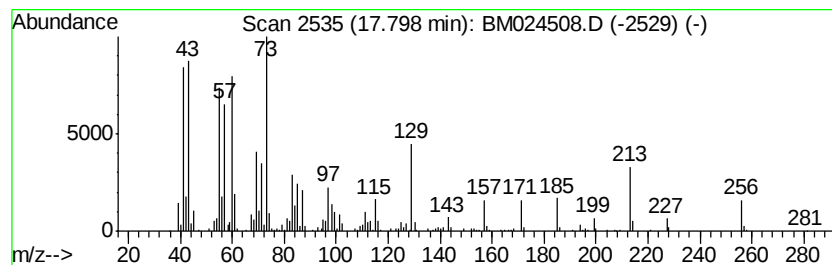
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.80	5.79 ng/ul	1020650	Phenanthrene-d10	16.85

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	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Tridecanoic acid	214	C13H26O2	000638-53-9	87
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

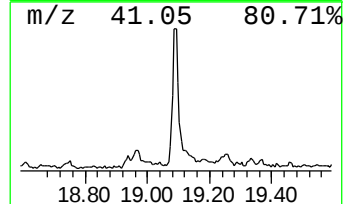
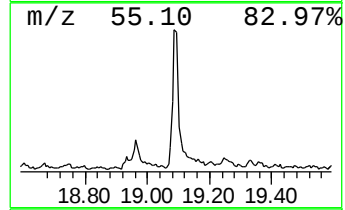
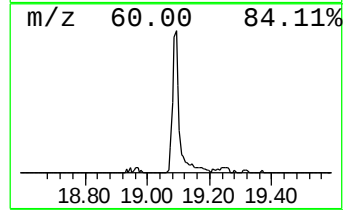
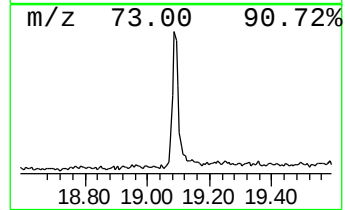
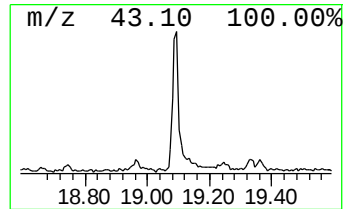
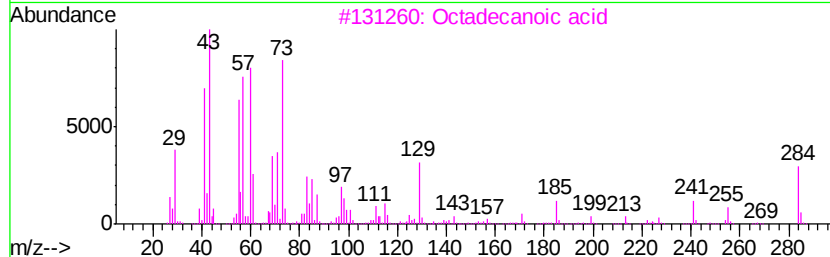
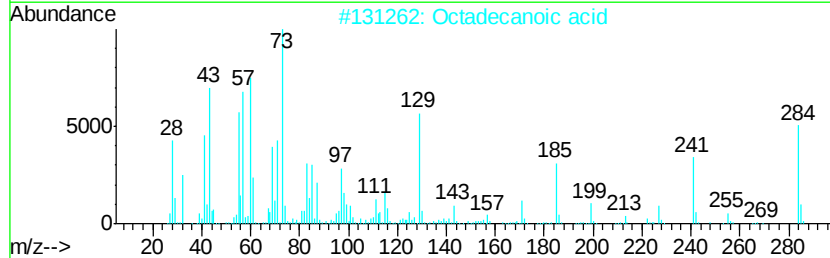
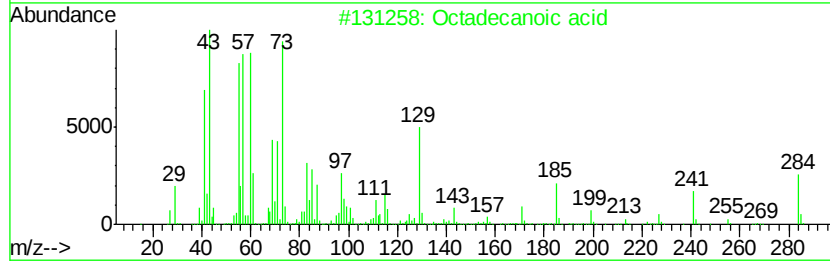
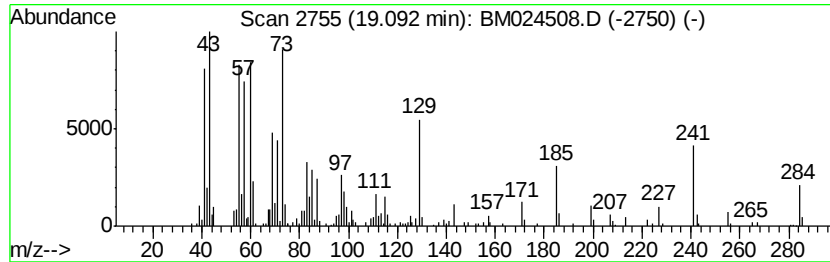
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	2.30 ng/ul	482156	Chrysene-d12	21.06

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	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

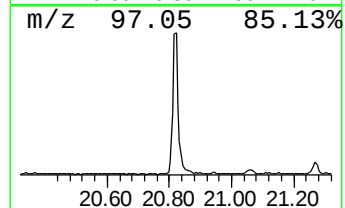
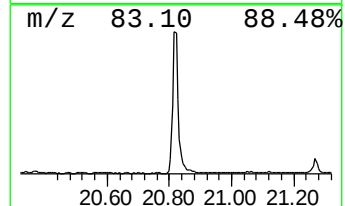
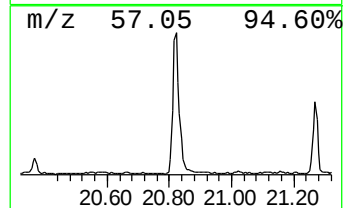
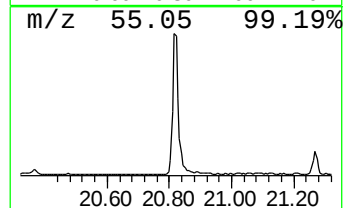
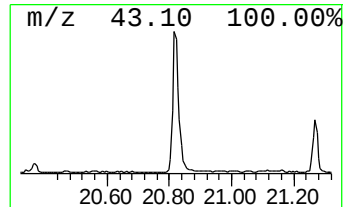
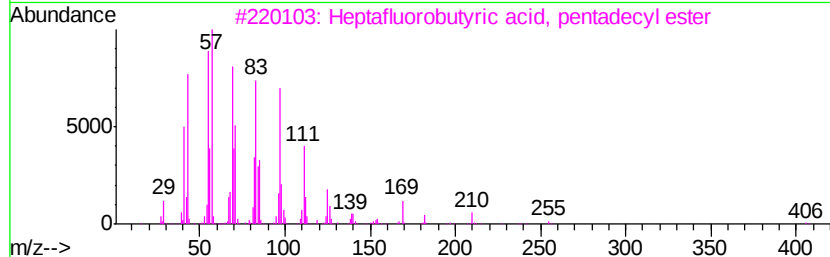
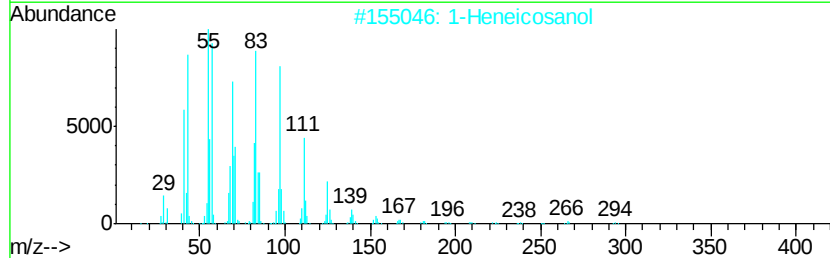
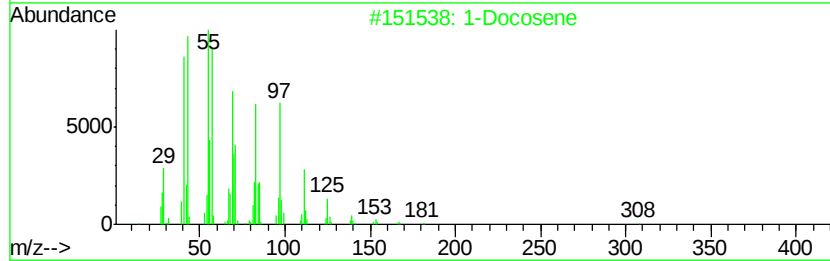
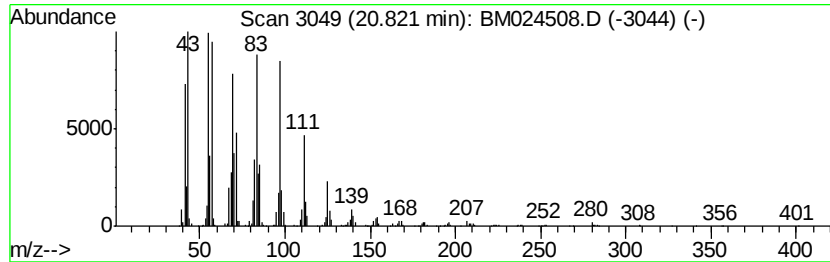
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic20.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	7.89 ng/ul	1654360	Chrysene-d12	21.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	99
2			1-Heneicosanol	312	C21H44O	015594-90-8	95
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

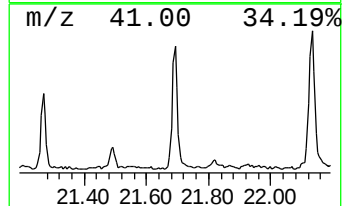
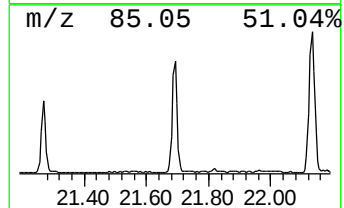
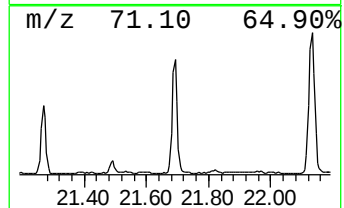
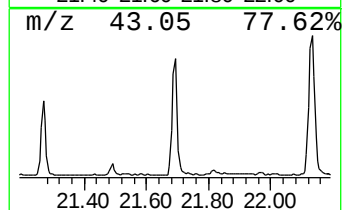
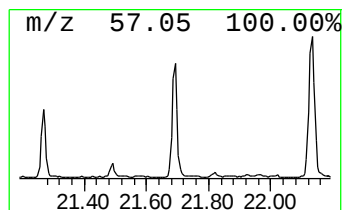
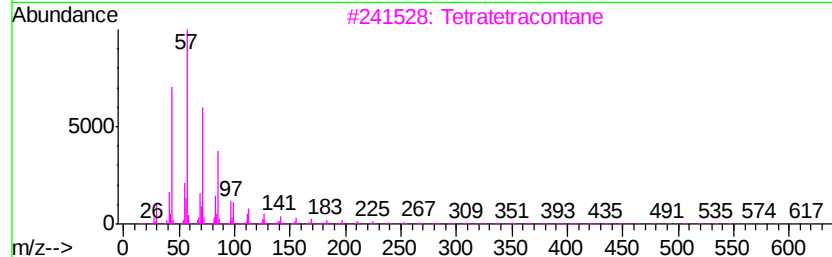
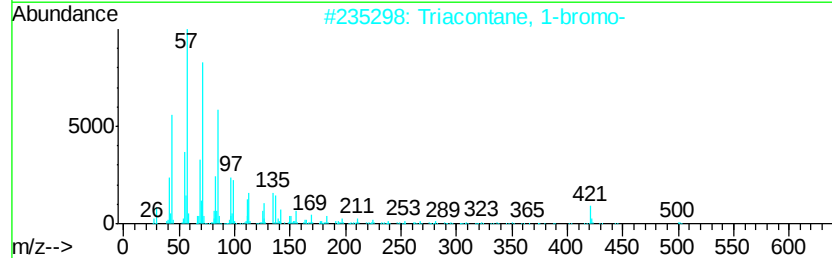
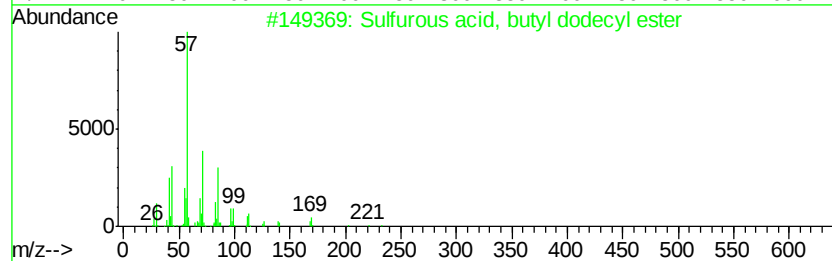
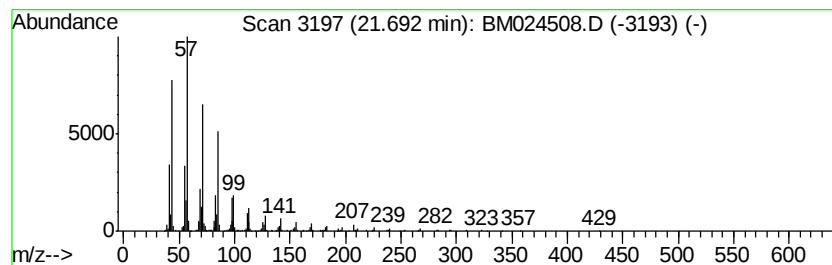
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Sulfurous acid, butyl dodec... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.69	2.87 ng/ul	602143	Chrysene-d12	21.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	91
2		Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Hentriacontane	437	C31H64	000630-04-6	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

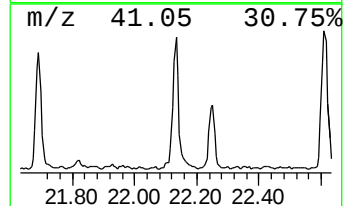
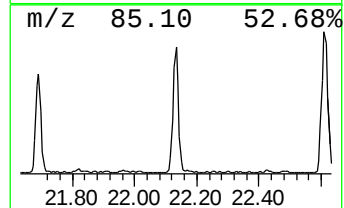
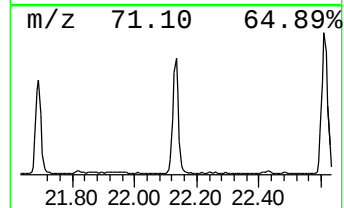
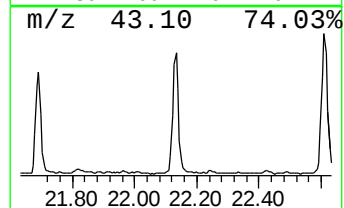
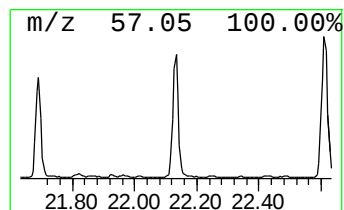
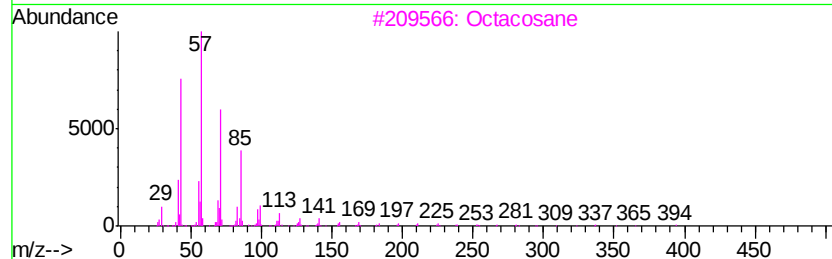
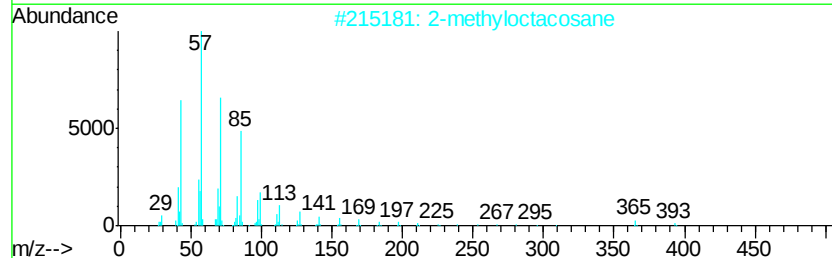
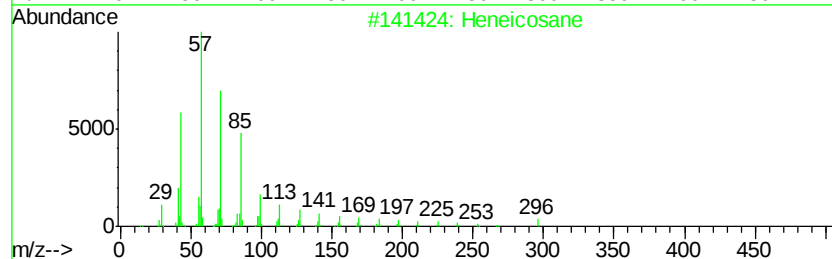
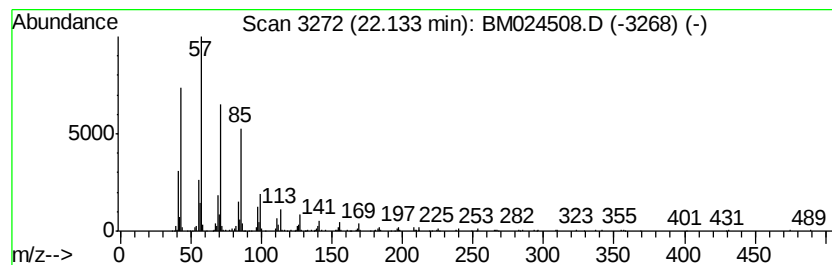
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.13	3.75 ng/ul	761222	Perylene-d12	23.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetracosane	338	C24H50	000646-31-1	90
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

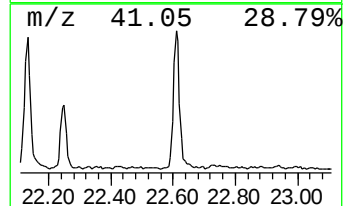
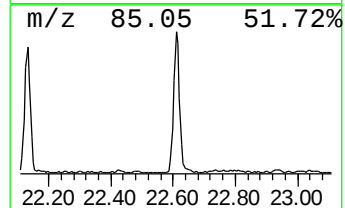
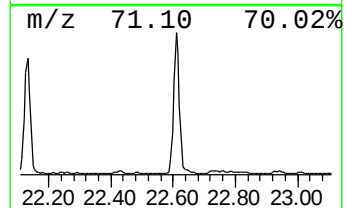
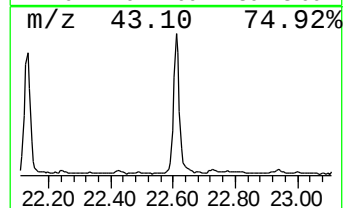
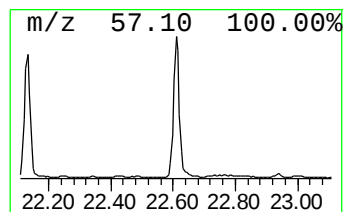
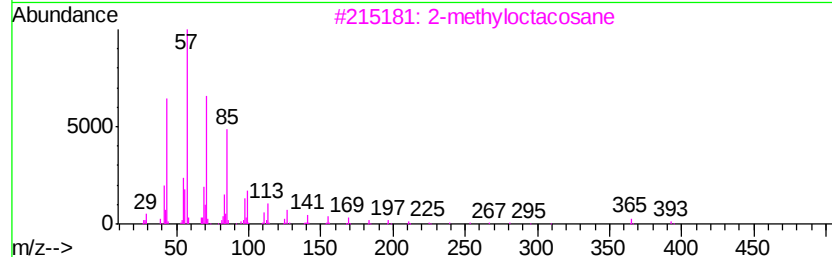
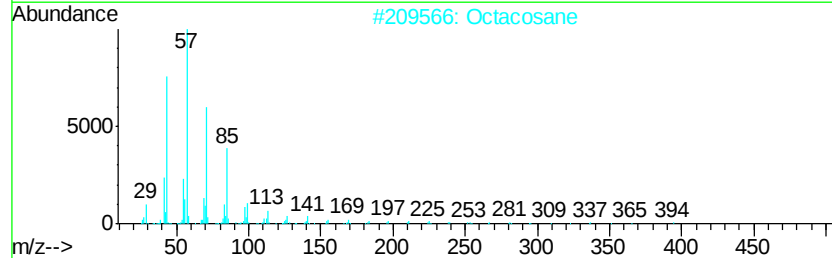
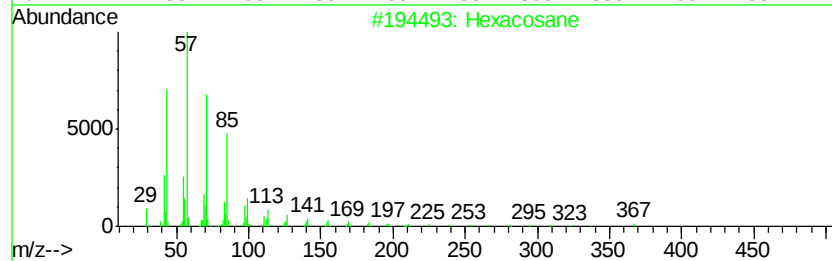
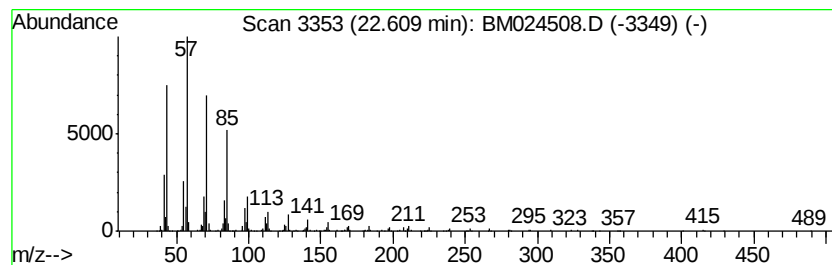
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.61	4.29 ng/ul	870388	Perylene-d12	23.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	93
2		Octacosane	394	C28H58	000630-02-4	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

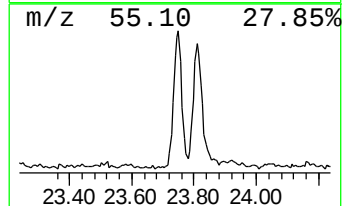
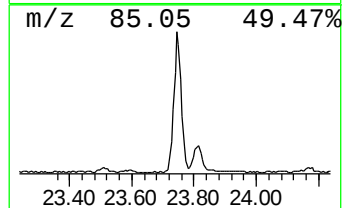
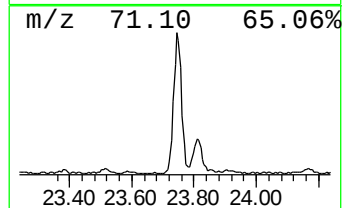
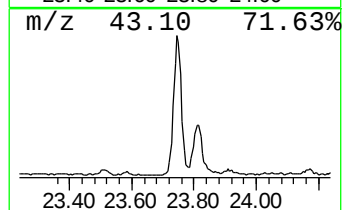
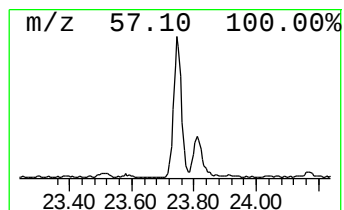
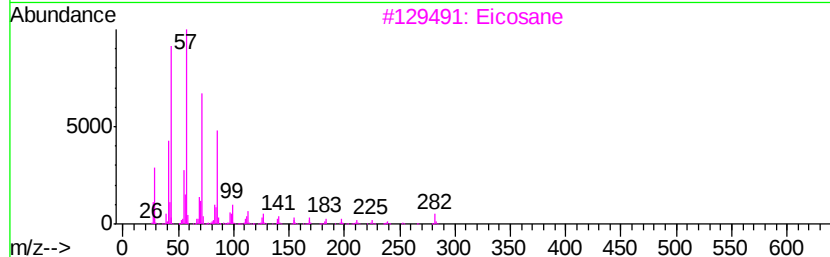
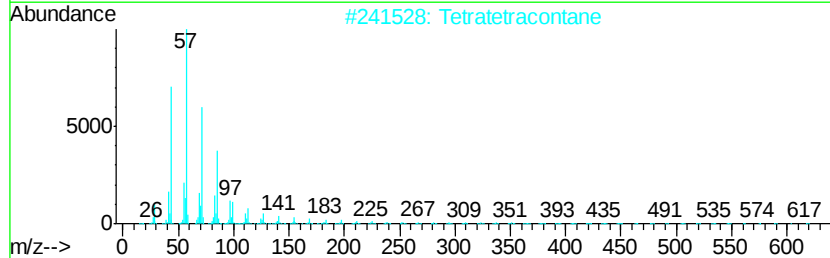
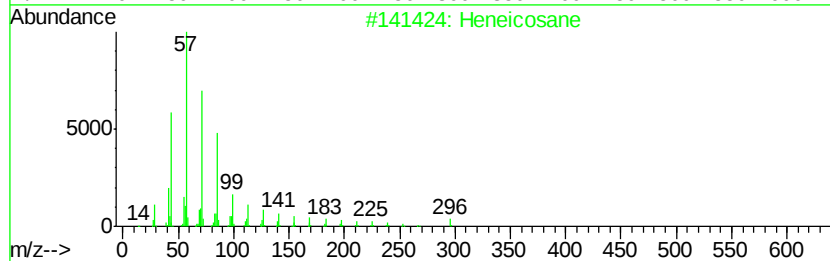
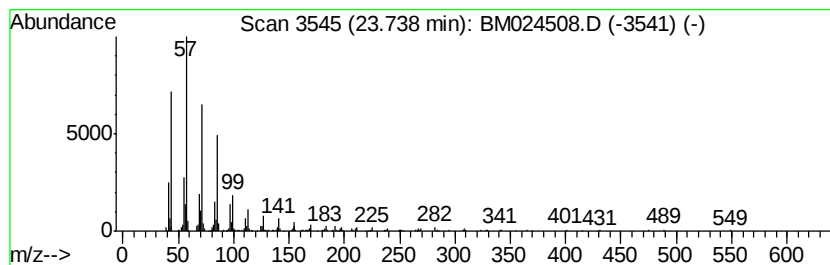
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
23.74	4.15 ng/ul	841064	Perylene-d12	23.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Tetratetracontane	619	C44H90	007098-22-8	93
3			Eicosane	282	C20H42	000112-95-8	92
4			Tetracosane	338	C24H50	000646-31-1	90
5			triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

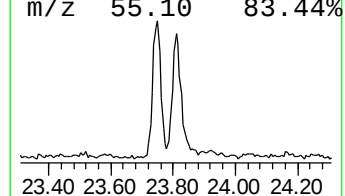
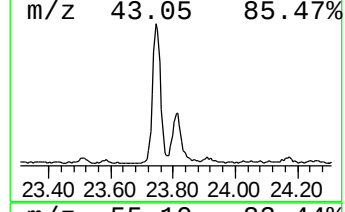
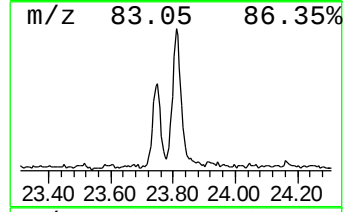
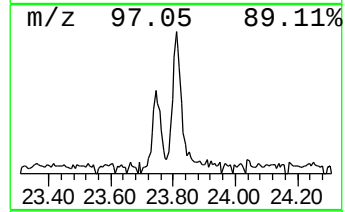
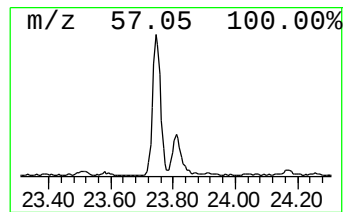
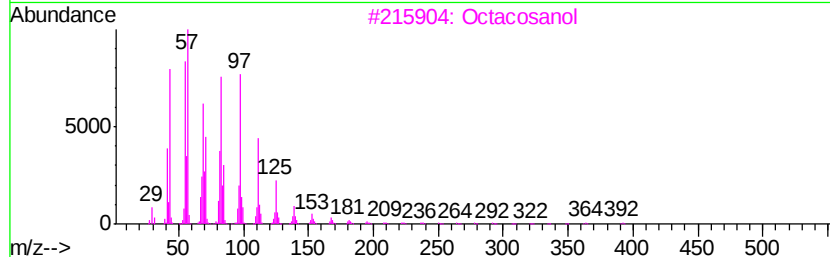
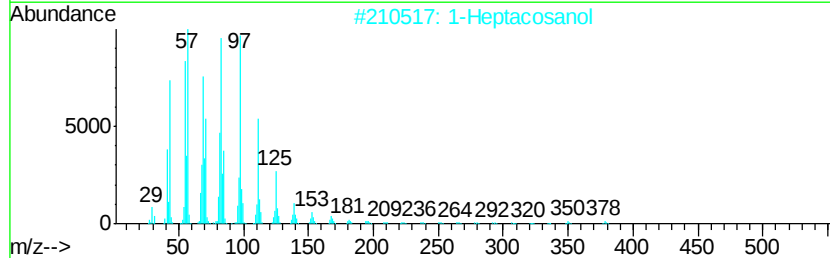
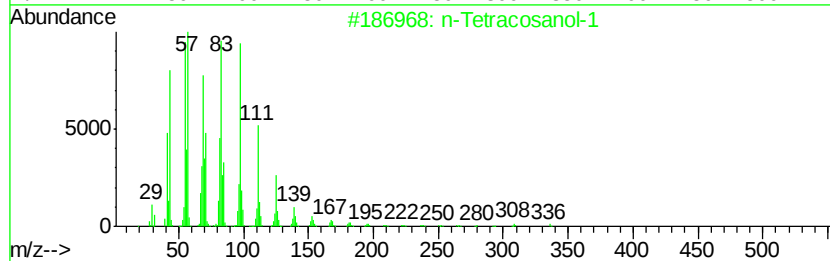
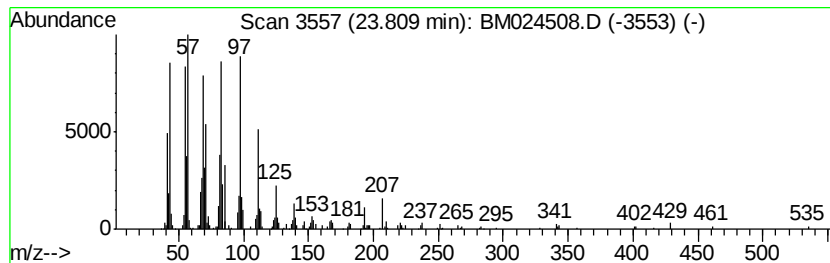
Instrument :
BNA_M
ClientSampleId :
GASG6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Tetracosanol-1 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.81	2.76 ng/ul	560315	Perylene-d12	23.18
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	n-Tetracosanol-1	354 C24H50O	000506-51-4	95
2	1-Heptacosanol	396 C27H56O	002004-39-9	95
3	Octacosanol	410 C28H58O	000557-61-9	93
4	1-Heneicosanol	312 C21H44O	015594-90-8	93
5	Behenic alcohol	326 C22H46O	000661-19-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
Data File : BM024508.D
Acq On : 17 Jan 2020 22:46
Operator : JU
Sample : L1109-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GASG6

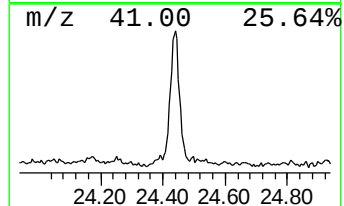
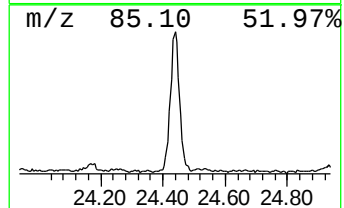
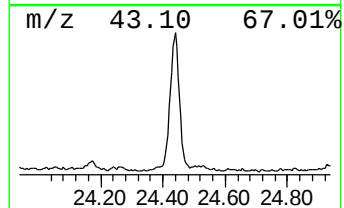
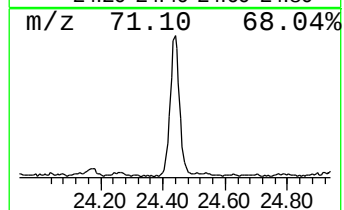
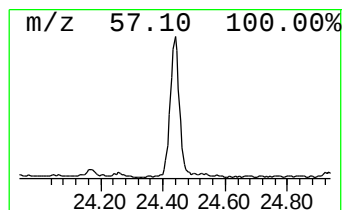
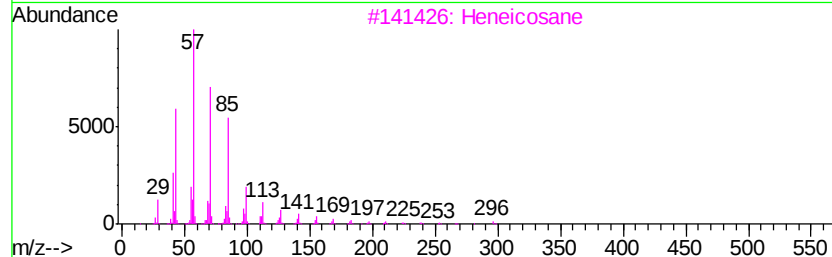
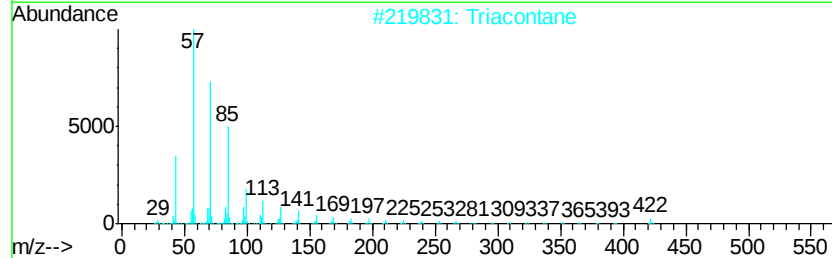
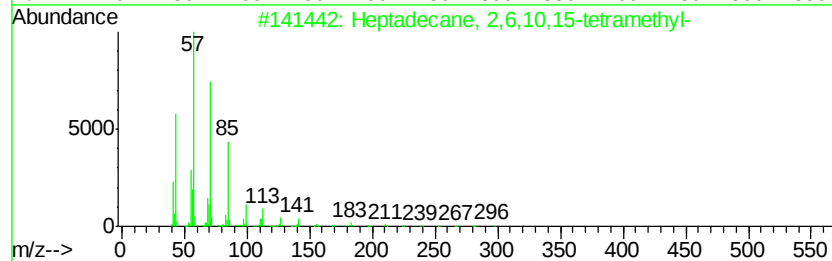
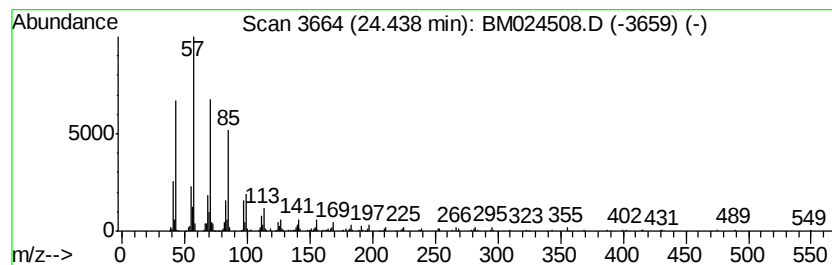
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.44	3.15 ng/ul	639339	Perylene-d12	23.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
2		Triacontane	422	C30H62	000638-68-6	90
3		Heneicosane	296	C21H44	000629-94-7	87
4		Tetratetracontane	619	C44H90	007098-22-8	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM011720\
 Data File : BM024508.D
 Acq On : 17 Jan 2020 22:46
 Operator : JU
 Sample : L1109-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GASG6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-ethyl-	7.55	2.4	ng/ul	144104	1	7.46	1211990	20.0
n-Hexadecanoic acid	17.80	5.8	ng/ul	1020650	4	16.85	3523850	20.0
Octadecanoic acid	19.09	2.3	ng/ul	482156	5	21.06	4191160	20.0
(DEL) Alkane: Cyc...	20.82	7.9	ng/ul	1654360	5	21.06	4191160	20.0
Sulfurous acid, b...	21.69	2.9	ng/ul	602143	5	21.06	4191160	20.0
(DEL) Alkane: Str...	22.13	3.8	ng/ul	761222	6	23.18	4056420	20.0
(DEL) Alkane: Str...	22.61	4.3	ng/ul	870388	6	23.18	4056420	20.0
(DEL) Alkane: Str...	23.74	4.2	ng/ul	841064	6	23.18	4056420	20.0
n-Tetracosanol-1	23.81	2.8	ng/ul	560315	6	23.18	4056420	20.0
(DEL) Alkane: Str...	24.44	3.1	ng/ul	639339	6	23.18	4056420	20.0