

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
 Data File : BN009848.D
 Acq On : 22 Feb 2020 16:34
 Operator : CG/JU
 Sample : L1516-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBD28

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_N\METHODS\SOM-EPA-BN021220MA.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	3.034	5	8	17	rVB	37784	53147	1.92%	0.143%
2	3.634	106	110	114	rBV7	2685	4541	0.16%	0.012%
3	3.716	122	124	129	rVB3	4826	6267	0.23%	0.017%
4	3.940	156	162	164	rBV7	5124	7728	0.28%	0.021%
5	4.599	271	274	280	rVB5	8857	12864	0.46%	0.035%
6	6.593	606	613	629	rBV	645939	1069831	38.62%	2.873%
7	6.746	634	639	655	rVB	465157	729615	26.34%	1.959%
8	6.934	664	671	686	rVB	626327	1026269	37.05%	2.756%
9	7.387	742	748	760	rVB	516873	821709	29.66%	2.206%
10	7.481	760	764	771	rBV6	6915	11948	0.43%	0.032%
11	8.104	864	870	887	rVB	719039	1183765	42.73%	3.178%
12	8.534	936	943	954	rBV	621478	1032657	37.28%	2.773%
13	8.710	971	973	975	rBV3	3829	3652	0.13%	0.010%
14	8.975	1013	1018	1023	rVB2	10257	15642	0.56%	0.042%
15	9.246	1057	1064	1071	rBV	510884	850039	30.69%	2.282%
16	9.775	1147	1154	1171	rBV	668409	1192275	43.04%	3.201%
17	10.134	1208	1215	1224	rBV	704321	1151613	41.57%	3.092%
18	10.293	1235	1242	1256	rBV	551972	1142635	41.25%	3.068%
19	11.757	1486	1491	1500	rVB3	6993	14017	0.51%	0.038%
20	12.640	1634	1641	1644	rBV5	1857	3066	0.11%	0.008%
21	12.869	1675	1680	1684	rBV4	3136	4793	0.17%	0.013%
22	12.945	1689	1693	1699	rVB5	7156	10992	0.40%	0.030%
23	13.340	1756	1760	1762	rBV3	2734	3205	0.12%	0.009%
24	13.457	1774	1780	1785	rBV	1159501	1658211	59.86%	4.452%
25	13.504	1785	1788	1798	rVB	99188	146607	5.29%	0.394%
26	13.716	1817	1824	1836	rBV	1395436	1956797	70.64%	5.254%
27	13.969	1862	1867	1869	rBV4	3396	4401	0.16%	0.012%
28	14.028	1871	1877	1887	rVB	1002241	1458119	52.64%	3.915%
29	14.275	1913	1919	1939	rBV	363892	763013	27.55%	2.049%
30	14.487	1949	1955	1959	rBV6	8645	13393	0.48%	0.036%
31	14.863	2016	2019	2021	rBV3	5052	4710	0.17%	0.013%
32	14.922	2024	2029	2032	rBV4	5469	7008	0.25%	0.019%
33	15.034	2042	2048	2058	rBV	1680209	2373303	85.68%	6.372%
34	15.192	2069	2075	2086	rBV	492734	804404	29.04%	2.160%

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022220\
 Data File : BN009848.D
 Acq On : 22 Feb 2020 16:34
 Operator : CG/JU
 Sample : L1516-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBD28

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_N\METHODS\SOM-EPA-BN021220MA.M
 Title : SVOA CALIBRATION

35	15.275	2086	2089	2094	rVV3	19985	27682	1.00%	0.074%
36	15.710	2159	2163	2166	rBV3	1701	3147	0.11%	0.008%
37	15.792	2173	2177	2179	rBV3	4304	6561	0.24%	0.018%
38	15.822	2180	2182	2188	rVB4	4950	6433	0.23%	0.017%
39	16.445	2284	2288	2290	rBV3	2403	3882	0.14%	0.010%
40	16.581	2308	2311	2316	rVB4	5473	7635	0.28%	0.021%
41	16.786	2340	2346	2355	rVV	1278672	1684788	60.82%	4.524%
42	16.886	2357	2363	2379	rVV	1784950	2465099	88.99%	6.619%
43	17.022	2382	2386	2387	rVV3	4799	6774	0.24%	0.018%
44	17.204	2413	2417	2424	rVB2	17494	23620	0.85%	0.063%
45	17.651	2489	2493	2499	rVB4	5986	10560	0.38%	0.028%
46	17.716	2499	2504	2513	rBV	190283	322110	11.63%	0.865%
47	17.780	2513	2515	2517	rVB2	7818	6393	0.23%	0.017%
48	17.904	2534	2536	2541	rVB5	4301	6248	0.23%	0.017%
49	17.945	2541	2543	2545	rBV3	2995	3592	0.13%	0.010%
50	17.992	2548	2551	2552	rBV2	6291	5744	0.21%	0.015%
51	18.116	2570	2572	2575	rVV4	3738	4205	0.15%	0.011%
52	18.145	2575	2577	2581	rVB4	9333	10769	0.39%	0.029%
53	18.398	2617	2620	2622	rBV2	3690	4384	0.16%	0.012%
54	18.522	2638	2641	2643	rBV4	4044	4227	0.15%	0.011%
55	18.586	2648	2652	2654	rBV4	9944	10670	0.39%	0.029%
56	18.657	2661	2664	2669	rVB6	5758	8724	0.31%	0.023%
57	18.827	2690	2693	2697	rBV	16145	19572	0.71%	0.053%
58	18.875	2697	2701	2702	rBV4	5411	5981	0.22%	0.016%
59	19.016	2720	2725	2732	rBV4	33420	57505	2.08%	0.154%
60	19.080	2733	2736	2741	rVB2	15245	20567	0.74%	0.055%
61	19.192	2749	2755	2762	rBV2	2094824	2770030	100.00%	7.438%
62	19.286	2768	2771	2776	rVB6	10688	11905	0.43%	0.032%
63	19.751	2847	2850	2854	rBV3	23214	26890	0.97%	0.072%
64	19.786	2854	2856	2860	rVB3	11834	13266	0.48%	0.036%
65	20.286	2935	2941	2944	rBV4	24167	33559	1.21%	0.090%
66	20.392	2957	2959	2963	rBV4	8439	11241	0.41%	0.030%
67	20.704	3006	3012	3015	rBV	44670	90014	3.25%	0.242%
68	20.745	3015	3019	3026	rVV2	1014608	1257946	45.41%	3.378%
69	20.804	3026	3029	3036	rVV	194775	273380	9.87%	0.734%
70	20.874	3039	3041	3043	rVB2	18836	15197	0.55%	0.041%
71	21.004	3058	3063	3072	rBV2	1408540	1793146	64.73%	4.815%

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

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_N\METHODS\SOM-EPA-BN021220MA.M
Title : SVOA CALIBRATION

72	21.192	3091	3095	3101	rBV	126756	142137	5.13%	0.382%
73	21.610	3162	3166	3173	rBV2	189883	273505	9.87%	0.734%
74	22.039	3236	3239	3244	rBV	202806	249000	8.99%	0.669%
75	22.157	3256	3259	3263	rVB	74730	76275	2.75%	0.205%
76	22.510	3315	3319	3323	rBV	239645	306421	11.06%	0.823%
77	22.974	3392	3398	3403	rBV	1614874	2618214	94.52%	7.030%
78	23.027	3403	3407	3413	rVV	209546	328868	11.87%	0.883%
79	23.104	3414	3420	3429	rVB	1058596	1801464	65.03%	4.837%
80	23.610	3502	3506	3512	rVB	175380	259902	9.38%	0.698%
81	23.680	3513	3518	3527	rVB2	230839	427578	15.44%	1.148%
82	24.274	3615	3619	3627	rVB2	102926	184465	6.66%	0.495%

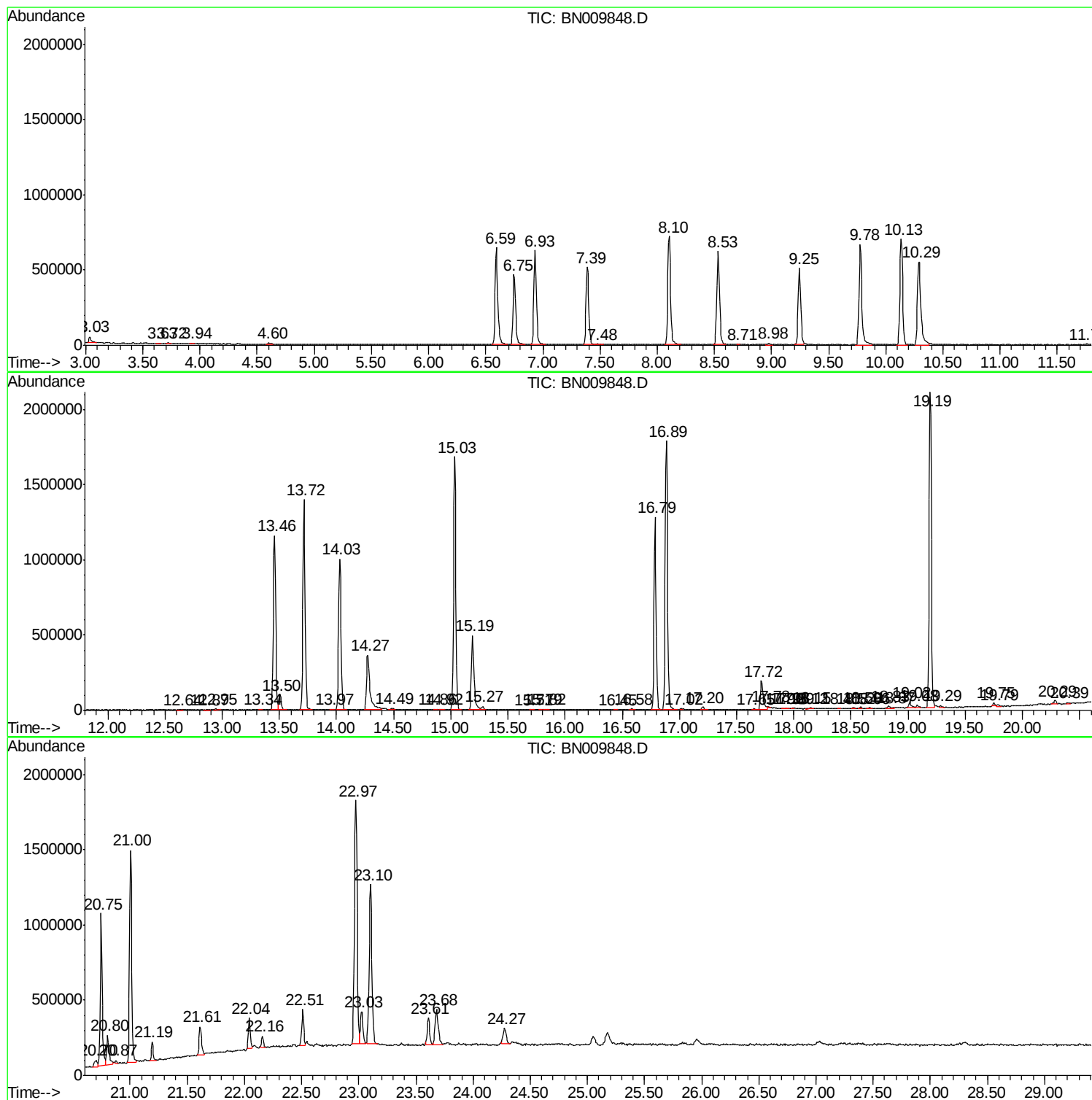
Sum of corrected areas: 37243511

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

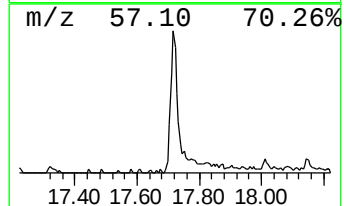
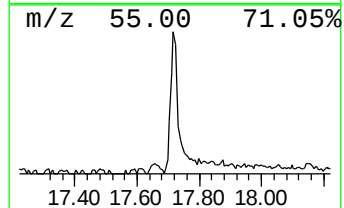
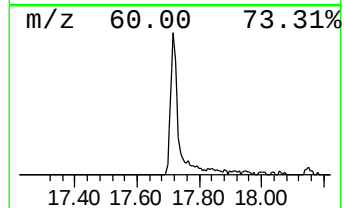
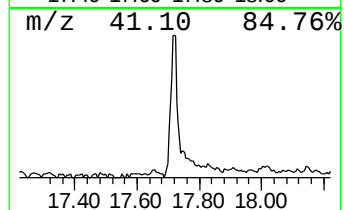
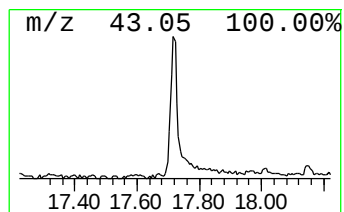
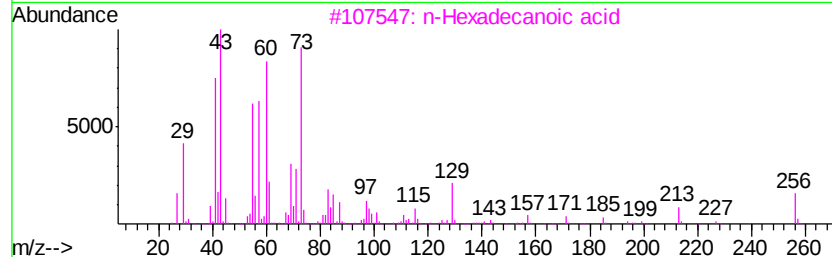
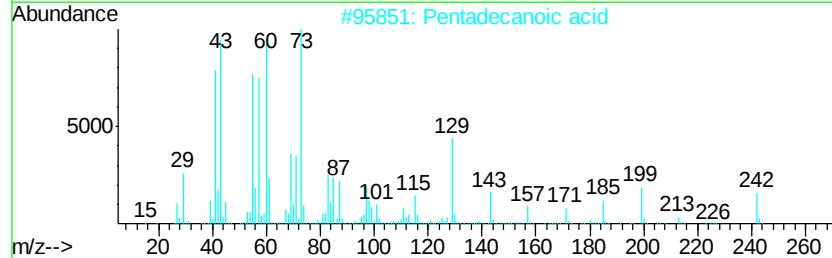
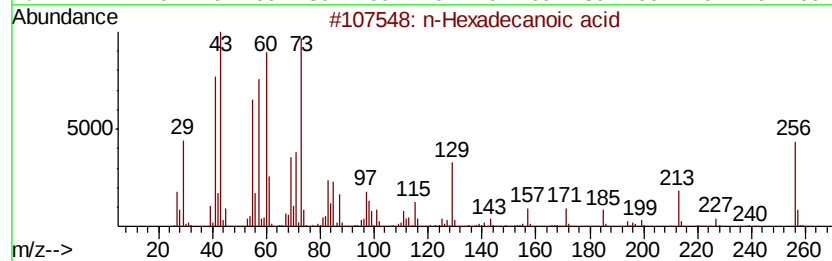
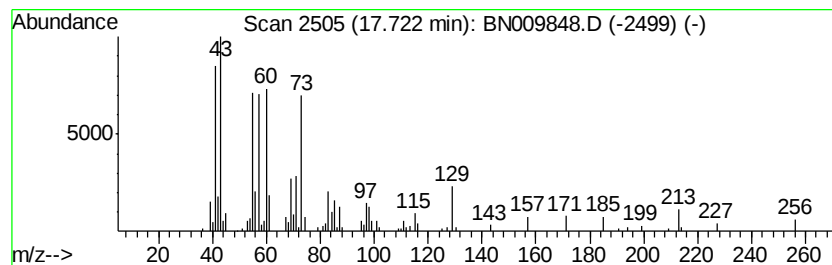
Instrument :
BNA_N
ClientSampleId :
DBD28

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.72	3.82 ng/ul	322110	Phenanthrene-d10		16.79	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5		Tridecanoic acid	214	C13H26O2	000638-53-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

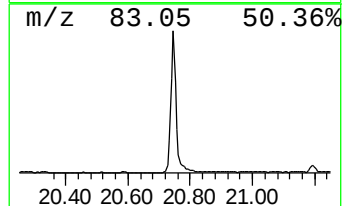
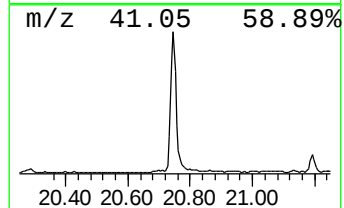
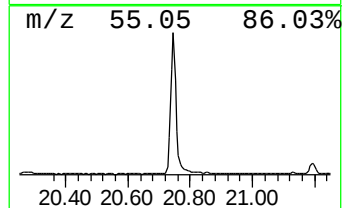
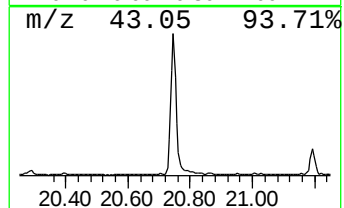
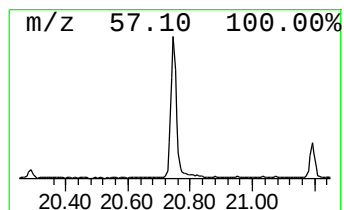
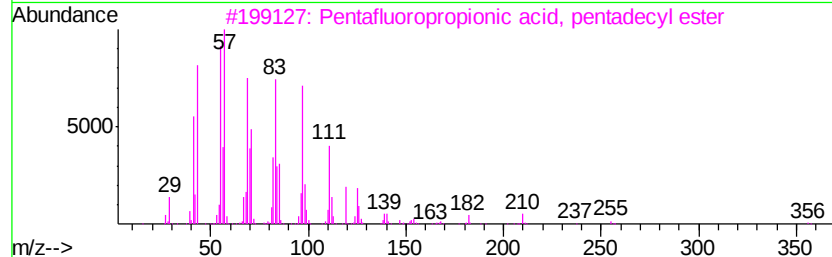
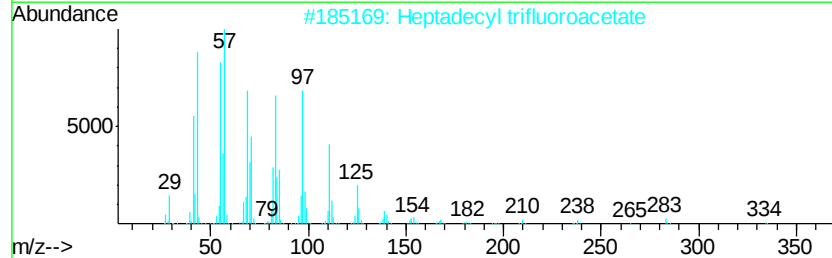
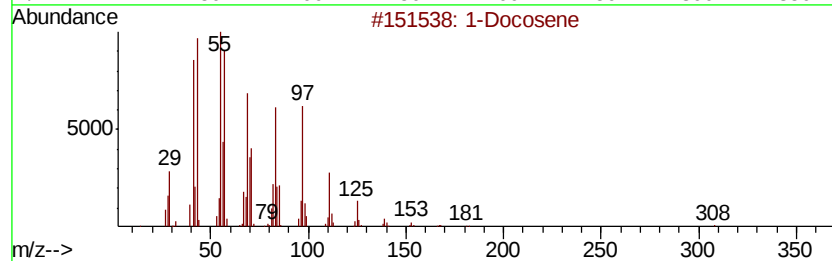
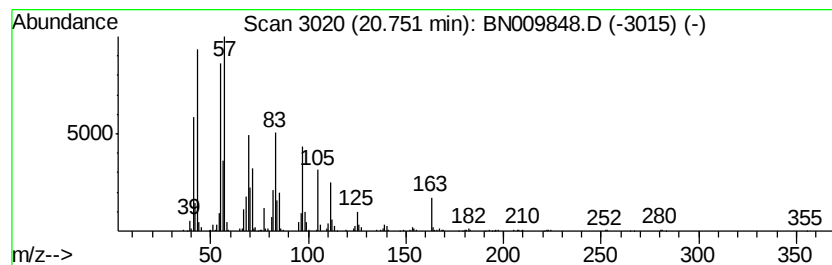
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	14.03 ng/ul	1257950	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	95
2		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
4		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

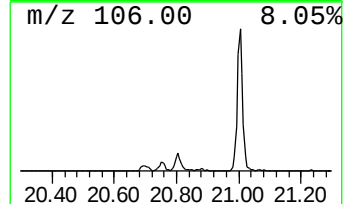
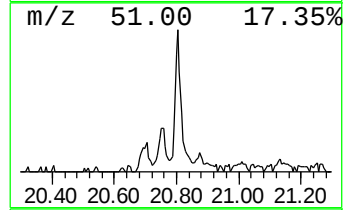
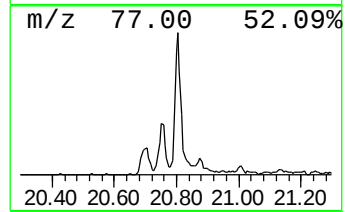
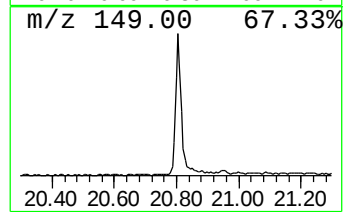
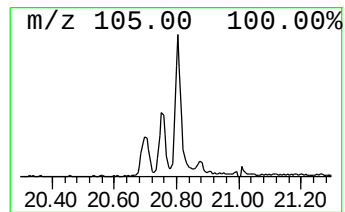
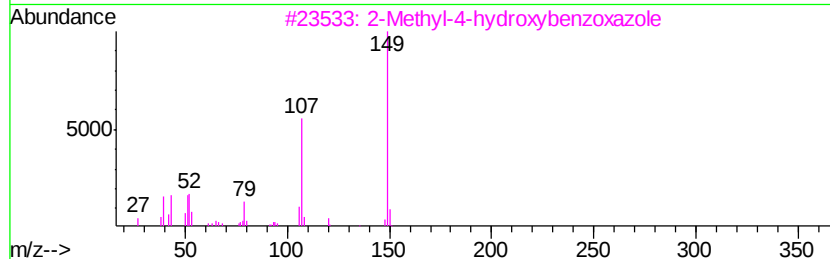
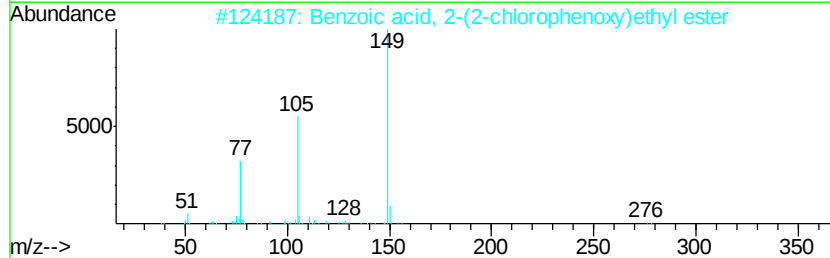
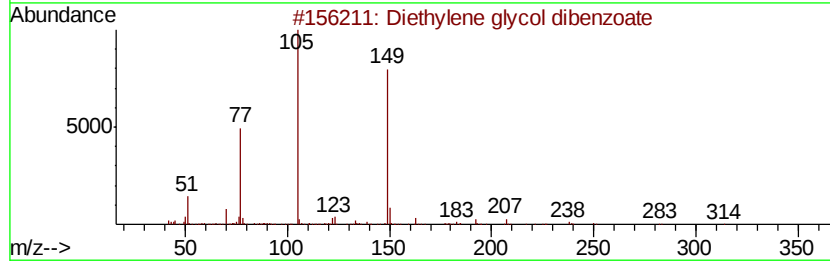
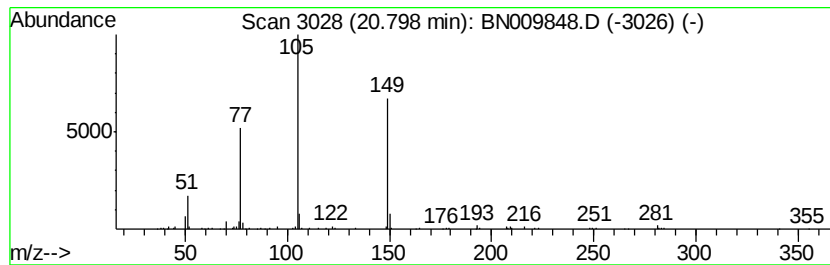
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Diethylene glycol dibenzoate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.80	3.05 ng/ul	273380	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	83
2		Benzoic acid, 2-(2-chlorophenoxy)ethyl ester	276	C15H13ClO3	1000368-25-9	78
3		2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	43
4		Vinyl benzoate	148	C9H8O2	000769-78-8	41
5		Cinnamic acid, .alpha.-[N-benzoyl-L-prolyl]	283	C16H13NO4	1000301-68-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

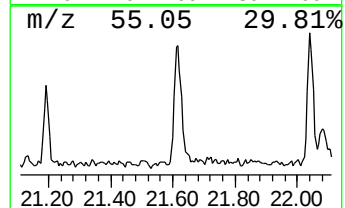
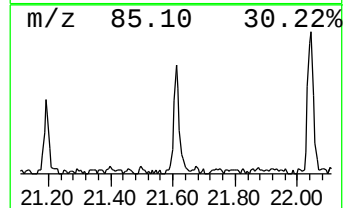
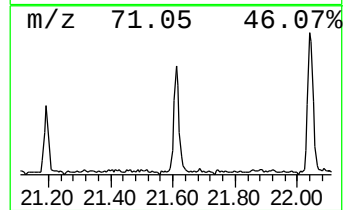
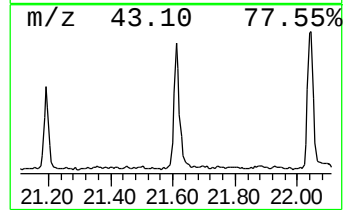
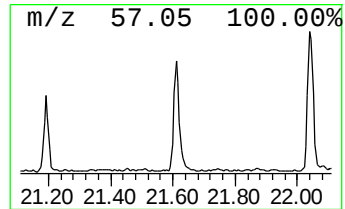
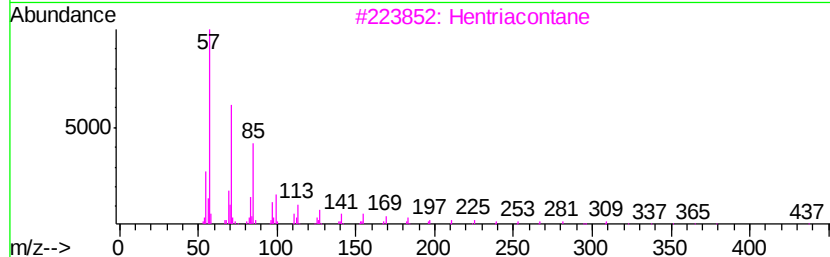
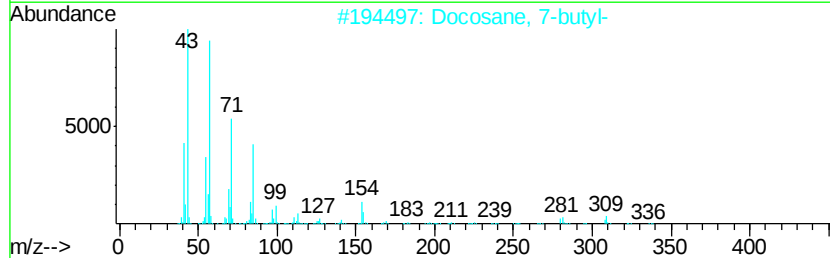
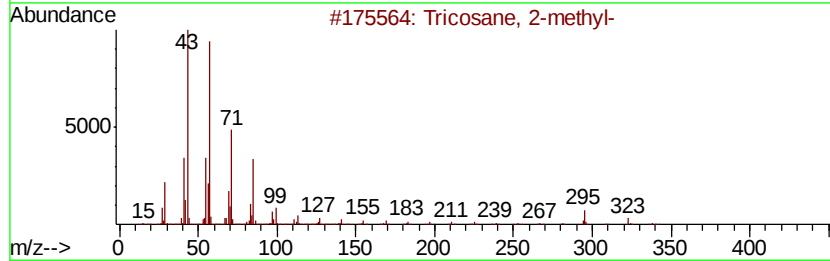
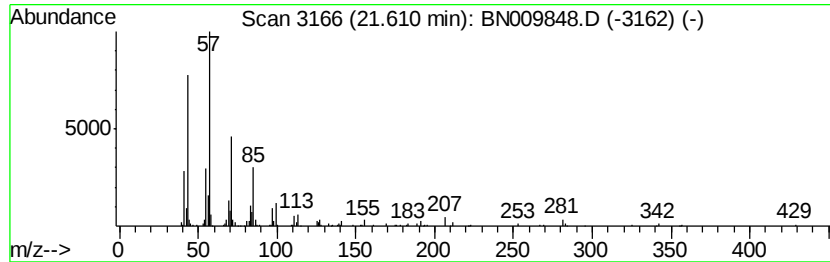
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.61	3.05 ng/ul	273505	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane, 2-methyl-	338	C24H50	001928-30-9	87
2		Docosane, 7-butyl-	366	C26H54	055282-15-0	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Heptacosane	380	C27H56	000593-49-7	86
5		Heptacosane	380	C27H56	000593-49-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

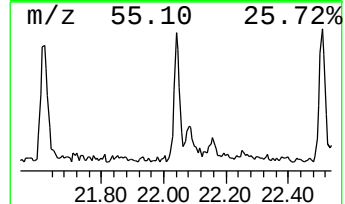
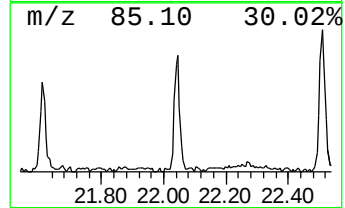
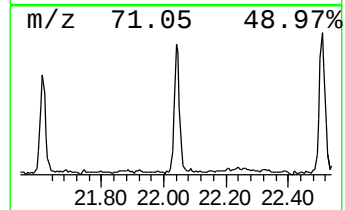
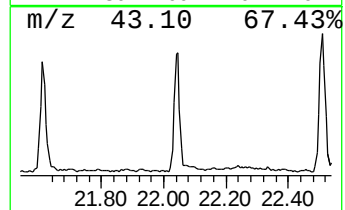
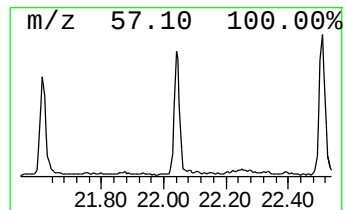
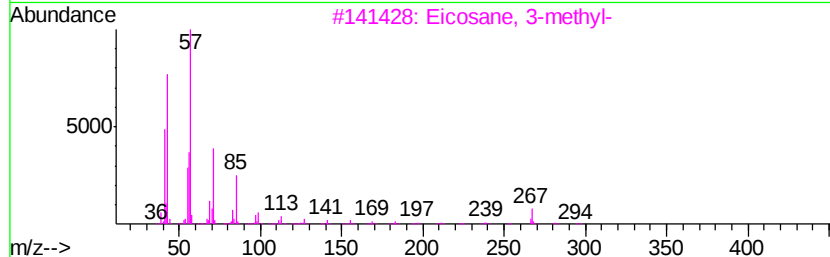
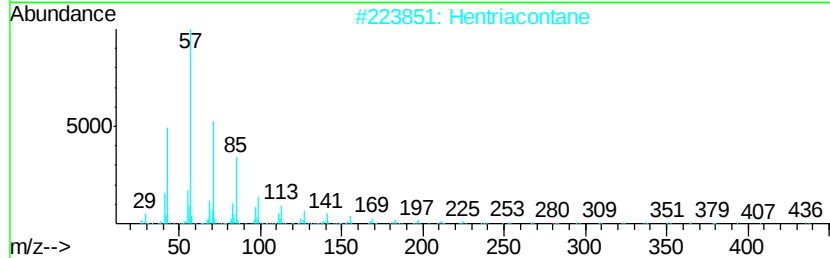
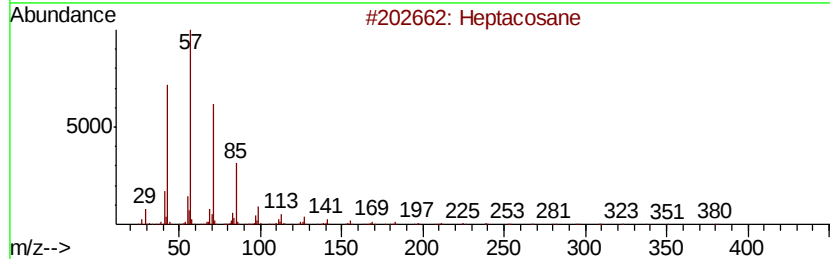
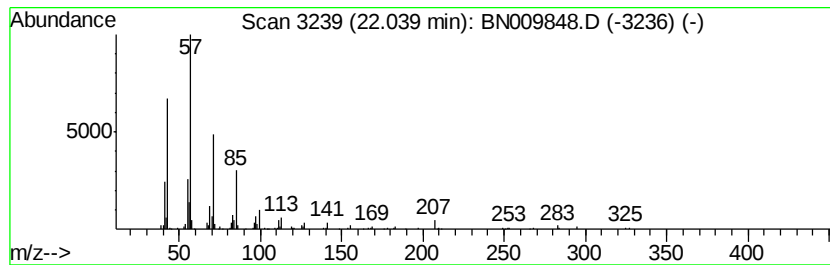
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.04	2.78 ng/ul	249000	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	90
2		Hentriacontane	437	C31H64	000630-04-6	86
3		Eicosane, 3-methyl-	296	C21H44	006418-46-8	86
4		Heptacosane	380	C27H56	000593-49-7	83
5		Hexatriacontane	507	C36H74	000630-06-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

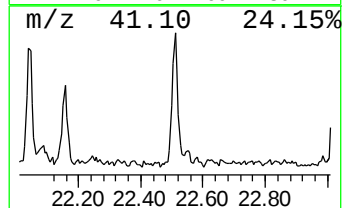
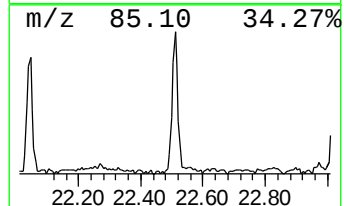
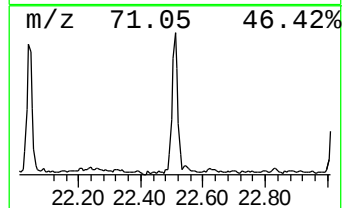
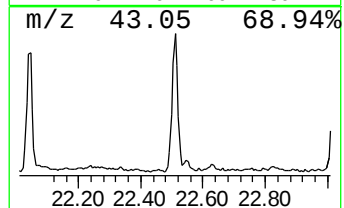
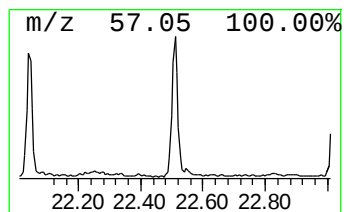
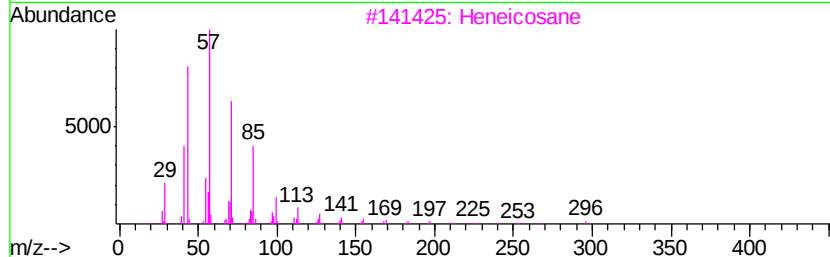
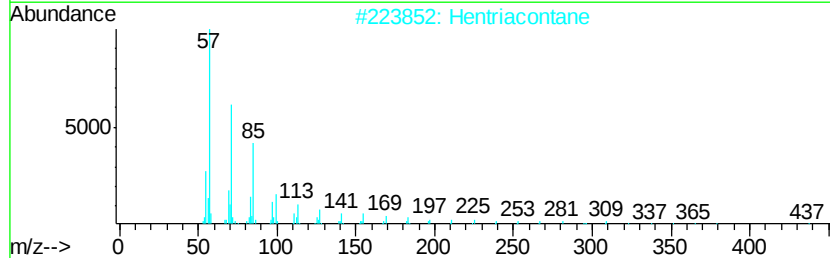
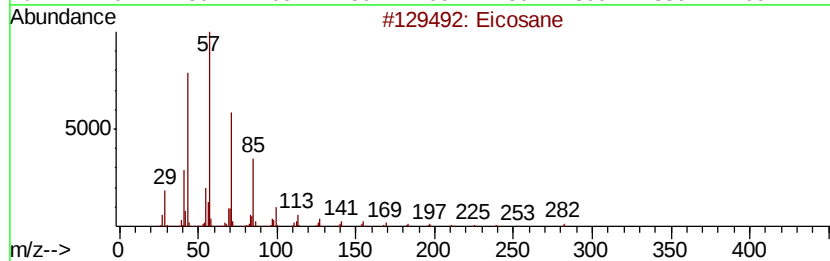
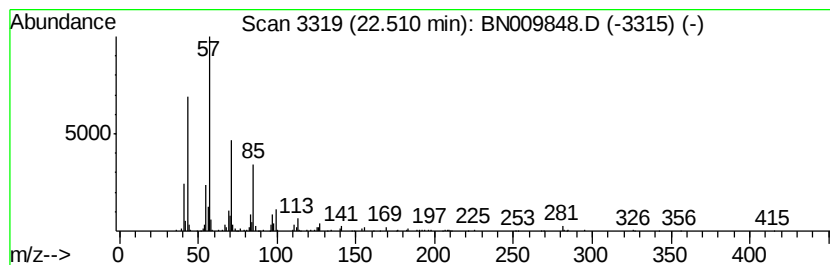
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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.
22.51	3.40 ng/ul	306421	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Hentriacontane	437	C31H64	000630-04-6	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Tetratetracontane	619	C44H90	007098-22-8	87
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

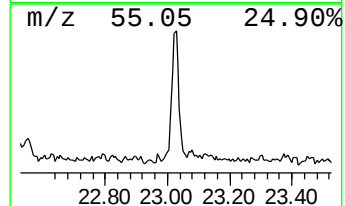
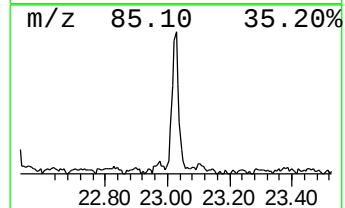
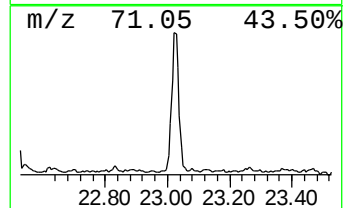
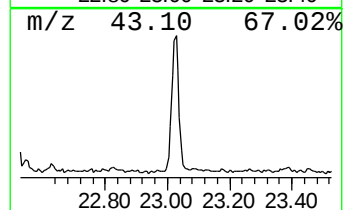
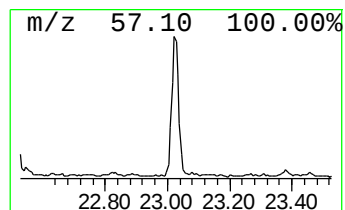
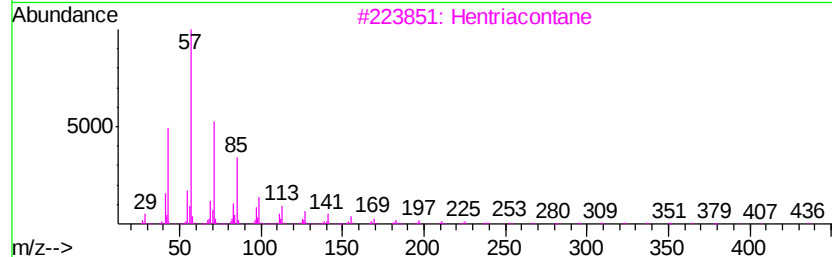
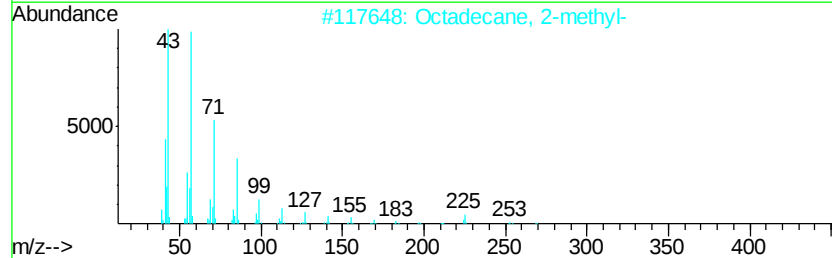
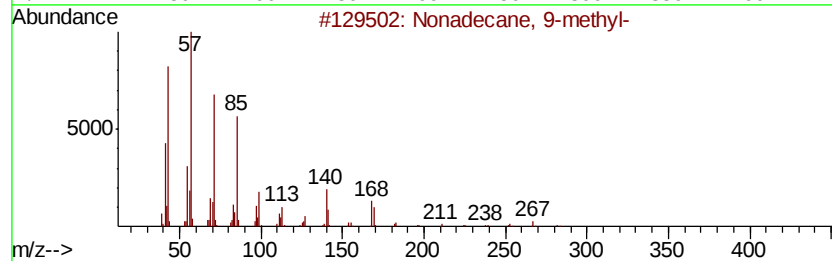
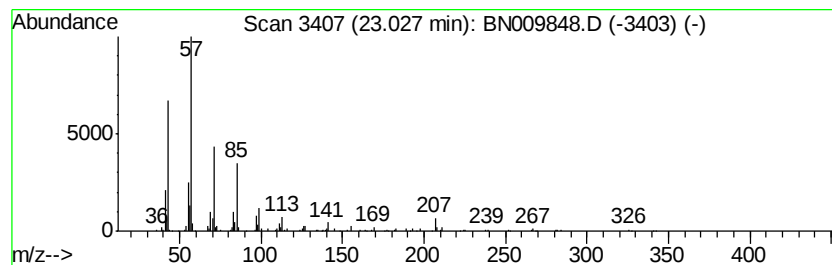
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.03	3.65 ng/ul	328868	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	87
3		Hentriacontane	437	C31H64	000630-04-6	86
4		Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	86
5		Tetratetracontane	619	C44H90	007098-22-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

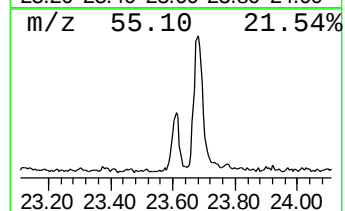
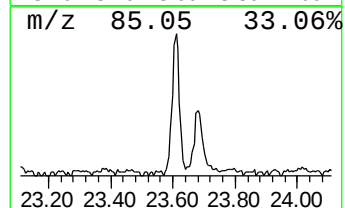
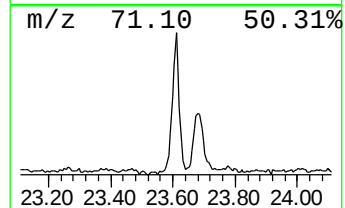
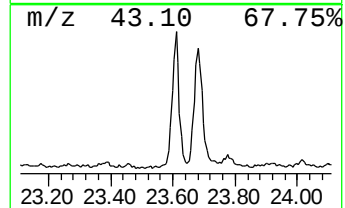
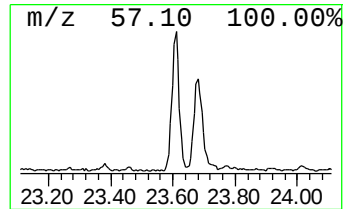
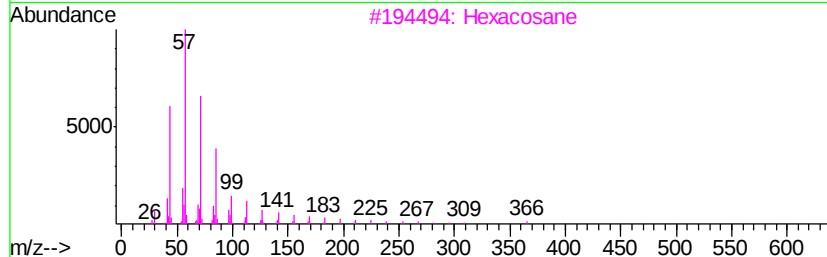
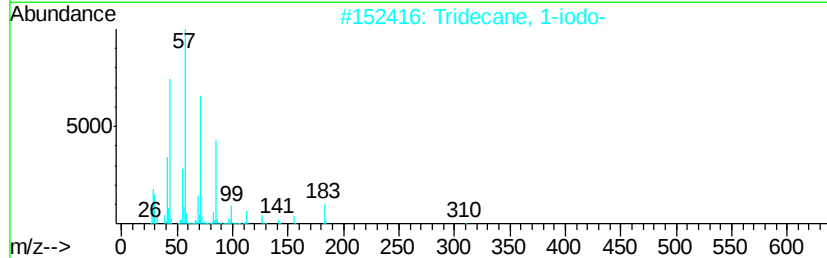
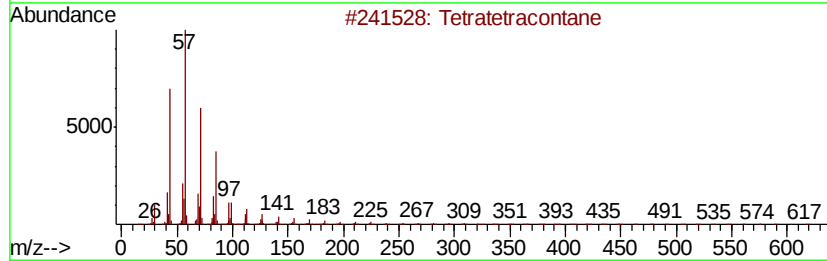
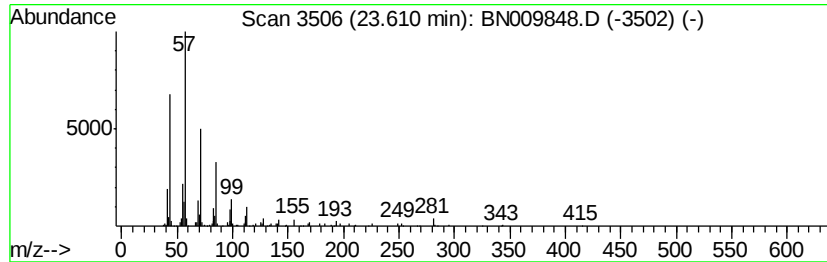
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	2.89 ng/ul	259902	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	86
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
3		Hexacosane	366	C26H54	000630-01-3	86
4		Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	83
5		2-methyloctacosane	408	C29H60	1000376-72-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

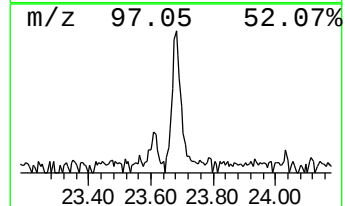
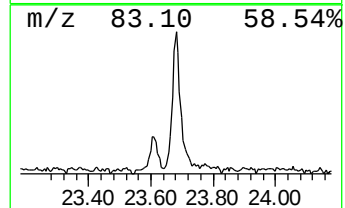
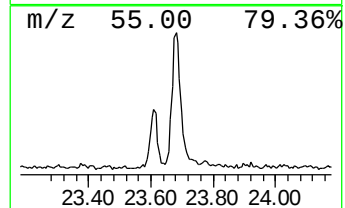
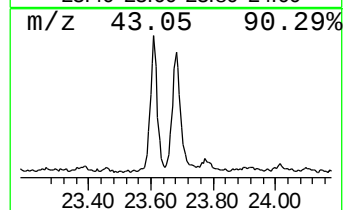
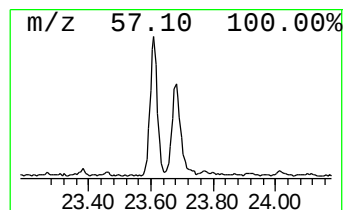
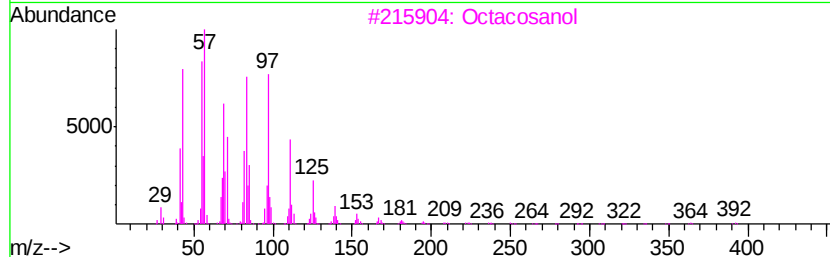
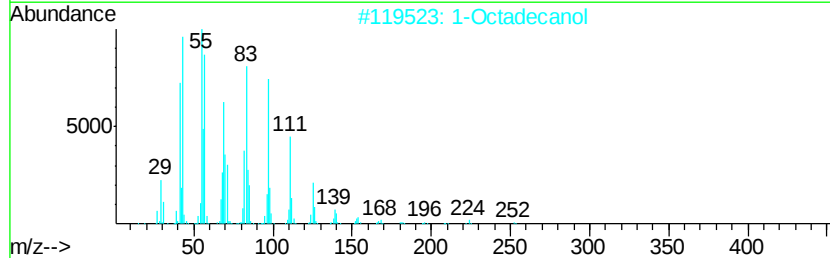
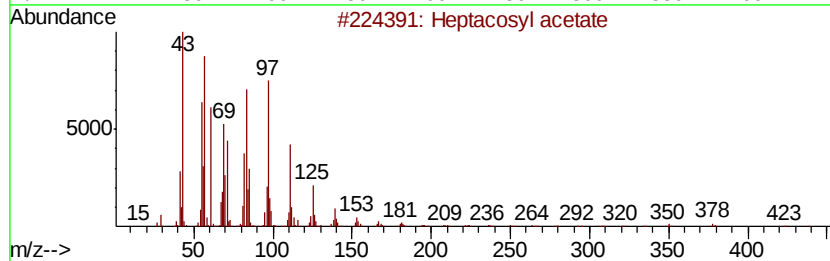
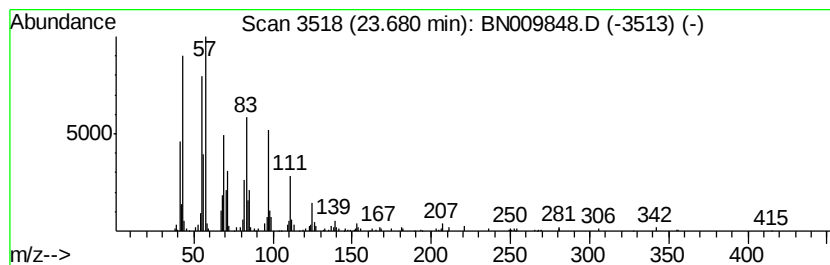
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Heptacosyl acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.68	4.75 ng/ul	427578	Perylene-d12	23.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosyl acetate	438	C29H58O2	1000351-78-2	87
2			1-Octadecanol	270	C18H38O	000112-92-5	87
3			Octacosanol	410	C28H58O	000557-61-9	87
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	87
5			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022220\
Data File : BN009848.D
Acq On : 22 Feb 2020 16:34
Operator : CG/JU
Sample : L1516-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD28

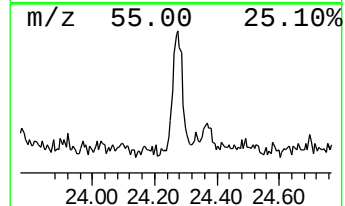
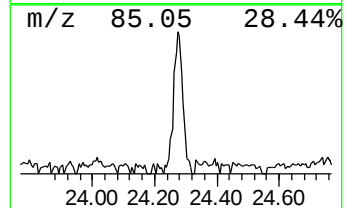
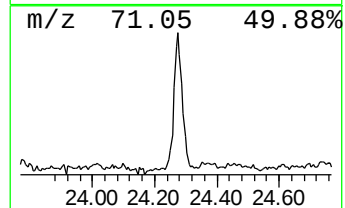
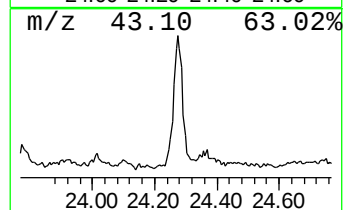
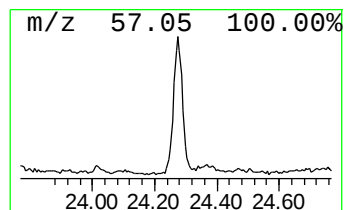
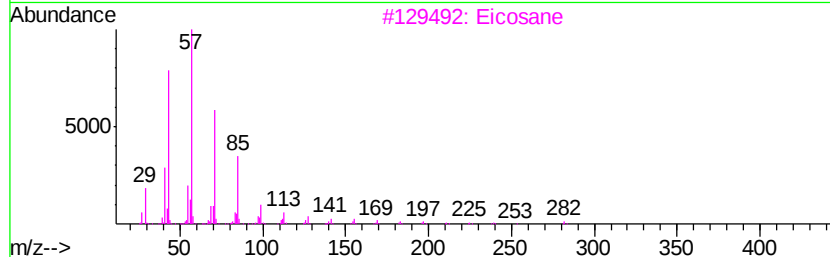
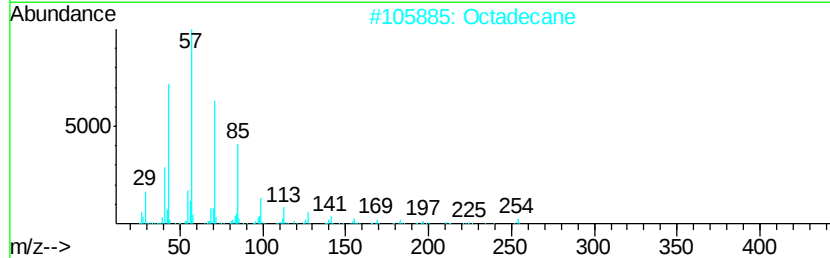
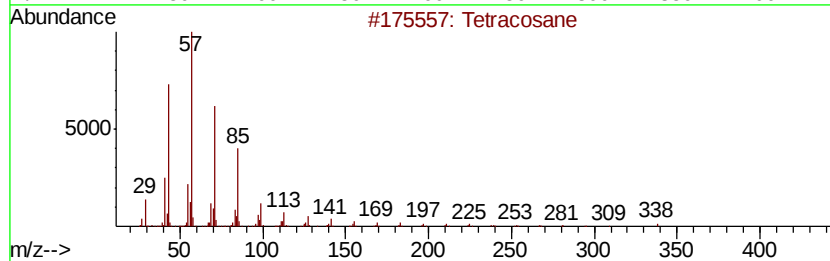
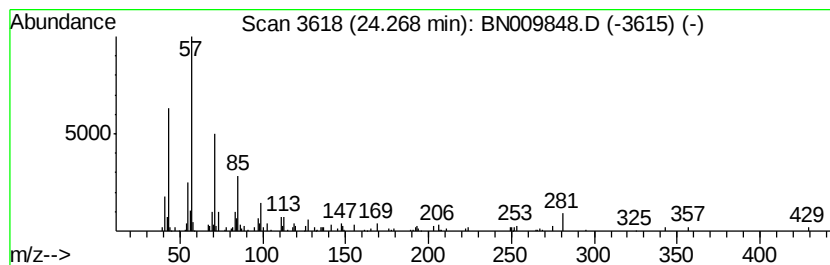
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.27	2.05 ng/ul	184465	Perylene-d12	23.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	81
2			Octadecane	254	C18H38	000593-45-3	74
3			Eicosane	282	C20H42	000112-95-8	74
4			Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	72
5			Nonadecane, 2-methyl-	282	C20H42	001560-86-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022220\
 Data File : BN009848.D
 Acq On : 22 Feb 2020 16:34
 Operator : CG/JU
 Sample : L1516-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBD28

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.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
n-Hexadecanoic acid	17.72	3.8	ng/ul	322110	4	16.79	1684790	20.0
(DEL) Alkane: Str...	20.75	14.0	ng/ul	1257950	5	21.00	1793150	20.0
Diethylene glycol...	20.80	3.0	ng/ul	273380	5	21.00	1793150	20.0
(DEL) Alkane: Str...	21.61	3.0	ng/ul	273505	5	21.00	1793150	20.0
(DEL) Alkane: Str...	22.04	2.8	ng/ul	249000	5	21.00	1793150	20.0
(DEL) Alkane: Str...	22.51	3.4	ng/ul	306421	6	23.10	1801460	20.0
(DEL) Alkane: Str...	23.03	3.6	ng/ul	328868	6	23.10	1801460	20.0
(DEL) Alkane: Str...	23.61	2.9	ng/ul	259902	6	23.10	1801460	20.0
Heptacosyl acetate	23.68	4.8	ng/ul	427578	6	23.10	1801460	20.0
(DEL) Alkane: Str...	24.27	2.0	ng/ul	184465	6	23.10	1801460	20.0