

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
 Data File : BM016154.D
 Acq On : 02 Aug 2018 00:16
 Operator : SJ/JU
 Sample : J4171-21
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JJ1L4

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 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-BM071318MA.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	7.351	685	691	710	rBV2	293733	715310	33.13%	3.072%
2	7.516	714	719	728	rBV	190748	296880	13.75%	1.275%
3	7.716	748	753	766	rBV	275041	460337	21.32%	1.977%
4	8.193	828	834	842	rBV	236787	383791	17.78%	1.648%
5	8.910	950	956	969	rBV	350090	594500	27.54%	2.553%
6	9.381	1030	1036	1049	rBV	289186	499520	23.14%	2.145%
7	10.098	1153	1158	1173	rBB	274324	469541	21.75%	2.016%
8	10.639	1244	1250	1266	rBV	367385	666003	30.85%	2.860%
9	11.010	1307	1313	1322	rBV	338282	577976	26.77%	2.482%
10	11.157	1332	1338	1359	rBV	334719	606250	28.08%	2.603%
11	14.227	1854	1860	1864	rBV	830445	1081288	50.09%	4.643%
12	14.274	1864	1868	1874	rVB	263756	364884	16.90%	1.567%
13	14.516	1903	1909	1920	rBV	799915	1195336	55.37%	5.133%
14	14.821	1954	1961	1969	rVB2	514752	787733	36.49%	3.383%
15	15.033	1992	1997	2018	rBV	291360	554012	25.66%	2.379%
16	15.810	2122	2129	2136	rBV	1114616	1519200	70.37%	6.524%
17	15.945	2147	2152	2165	rBV	377147	544845	25.24%	2.340%
18	17.551	2420	2425	2430	rBV	656461	921661	42.69%	3.958%
19	17.592	2430	2432	2436	rVV	39370	46842	2.17%	0.201%
20	17.651	2436	2442	2454	rVB	1269184	1709464	79.18%	7.341%
21	18.262	2541	2546	2556	rBV	235951	296205	13.72%	1.272%
22	18.398	2564	2569	2578	rBV	266899	382053	17.70%	1.641%
23	18.480	2578	2583	2589	rVB2	23707	38807	1.80%	0.167%
24	19.509	2754	2758	2761	rBV4	22245	33949	1.57%	0.146%
25	19.586	2767	2771	2775	rVB	53494	67474	3.13%	0.290%
26	19.656	2778	2783	2794	rBV	164934	245074	11.35%	1.052%
27	19.868	2815	2819	2821	rBV	27353	27583	1.28%	0.118%
28	19.915	2821	2827	2837	rVB	1557432	2158904	100.00%	9.271%
29	20.392	2904	2908	2912	rVB	56382	64164	2.97%	0.276%
30	20.803	2973	2978	2987	rVB4	110602	190997	8.85%	0.820%
31	20.880	2987	2991	2995	rBV	85847	94746	4.39%	0.407%
32	21.345	3065	3070	3083	rBV3	262511	539113	24.97%	2.315%
33	21.674	3120	3126	3130	rBV	865655	1144016	52.99%	4.913%
34	21.768	3139	3142	3147	rVB	130400	132714	6.15%	0.570%

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35	22.215	3215	3218	3222	rBV	120044	155486	7.20%	0.668%
36	22.256	3222	3225	3232	rVB3	63551	102550	4.75%	0.440%
37	22.697	3296	3300	3305	rBV	122118	164818	7.63%	0.708%
38	22.833	3321	3323	3328	rVB2	66121	83731	3.88%	0.360%
39	23.221	3385	3389	3396	rBV	118036	177837	8.24%	0.764%
40	23.815	3486	3490	3494	rBV3	94967	160505	7.43%	0.689%
41	23.880	3494	3501	3516	rVB	1007565	1925216	89.18%	8.267%
42	24.027	3519	3526	3535	rBV2	472902	962801	44.60%	4.135%
43	24.491	3601	3605	3611	rVB2	84111	142798	6.61%	0.613%

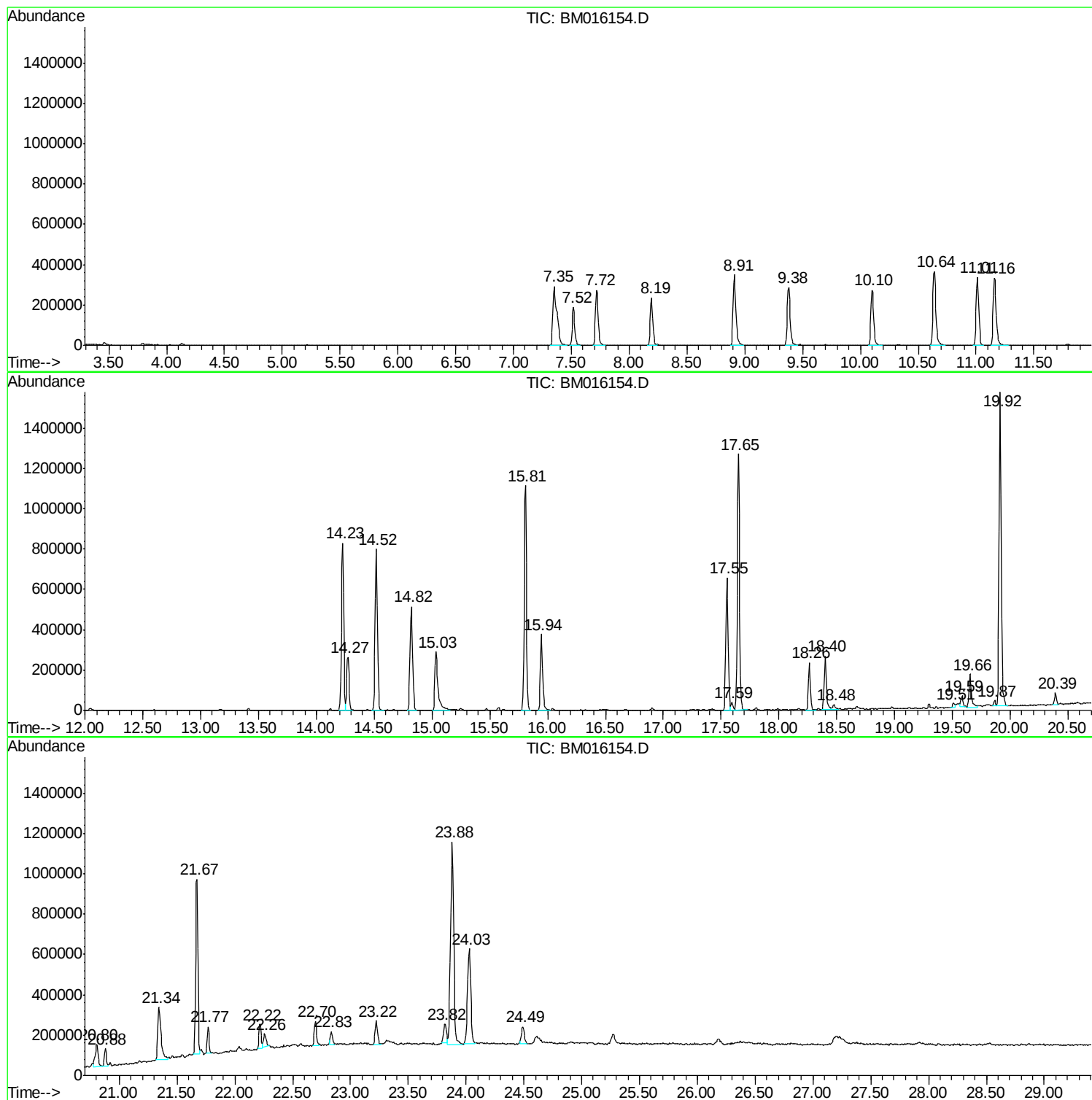
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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



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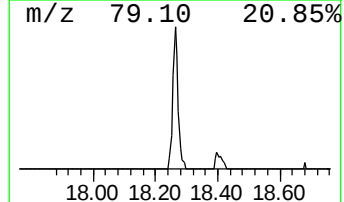
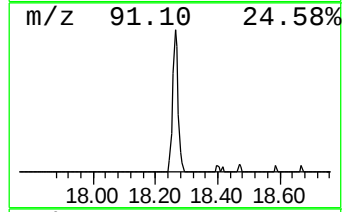
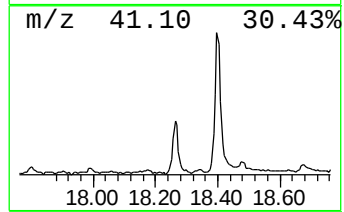
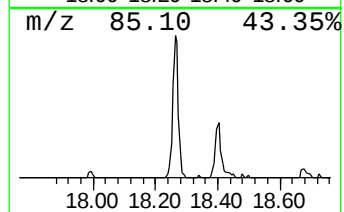
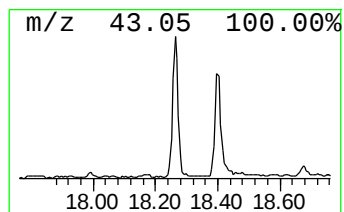
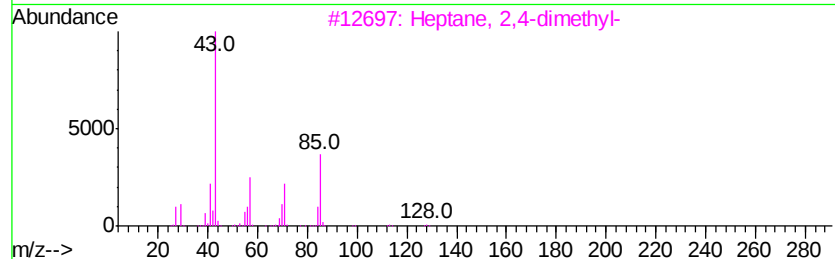
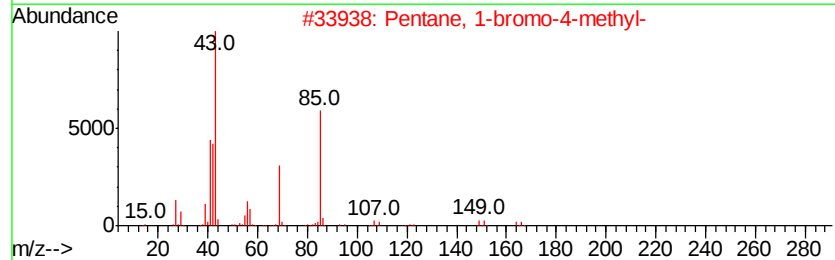
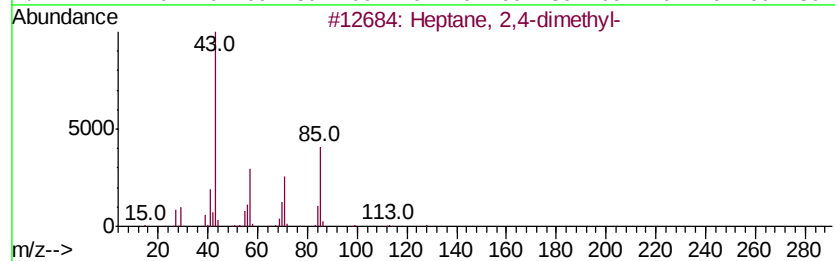
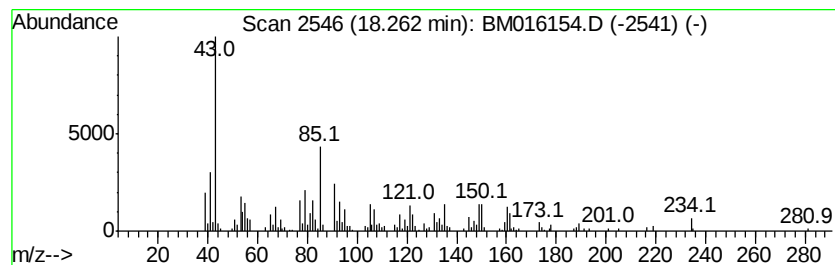
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.26	6.43 ng/ul	296205	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	14
2		Pentane, 1-bromo-4-methyl-	164	C6H13Br	000626-88-0	14
3		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	14
4		Hexane, 3-bromo-	164	C6H13Br	003377-87-5	14
5		Pentanoic acid, 2,2-dimethyl-, e...	156	C9H16O2	044970-05-0	14



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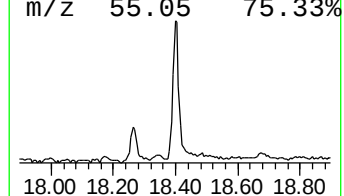
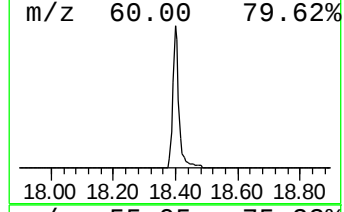
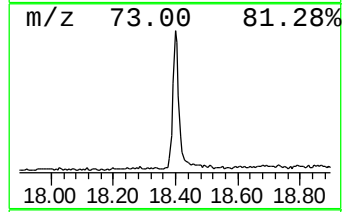
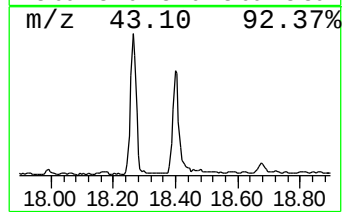
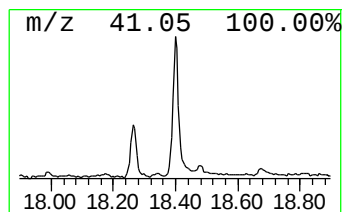
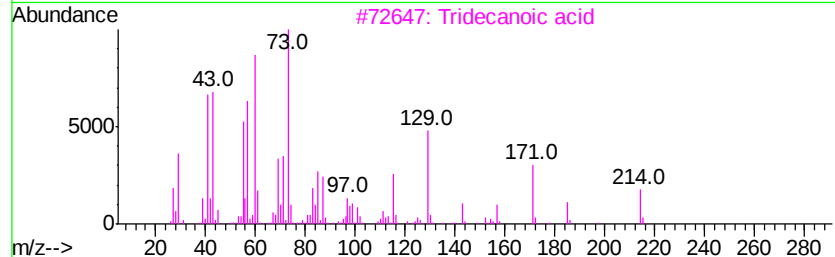
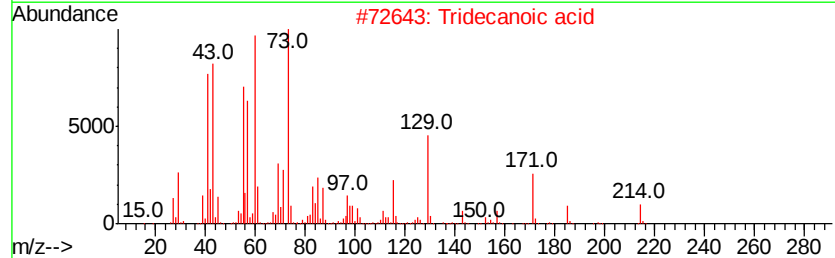
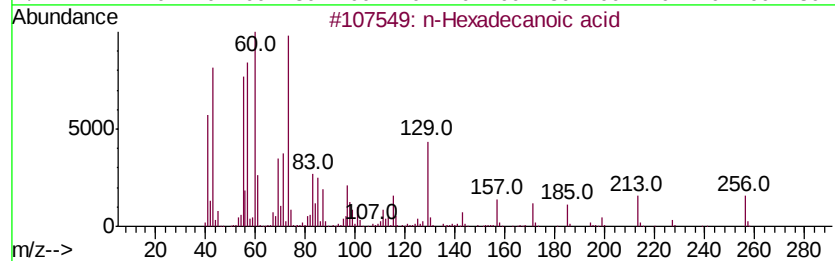
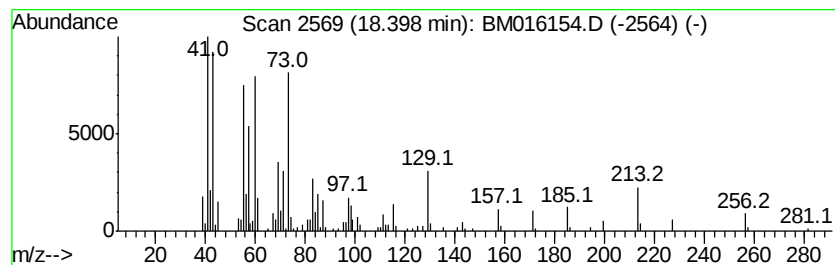
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.40	8.29 ng/ul	382053	Phenanthrene-d10		17.55	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Tridecanoic acid	214	C13H26O2	000638-53-9	89
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	74



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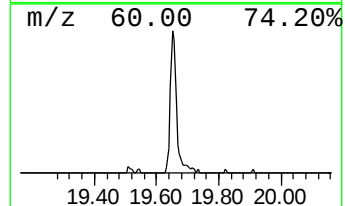
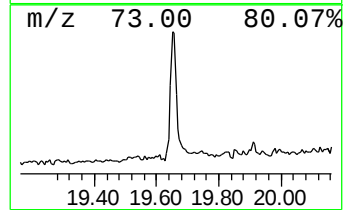
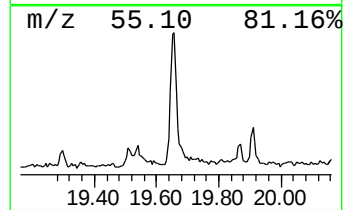
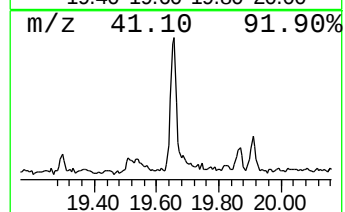
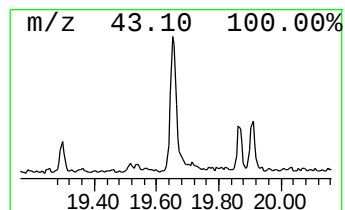
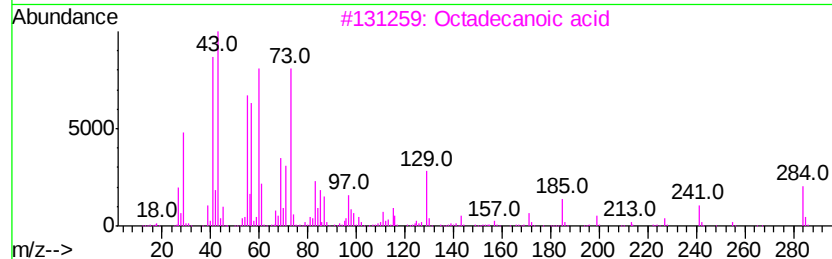
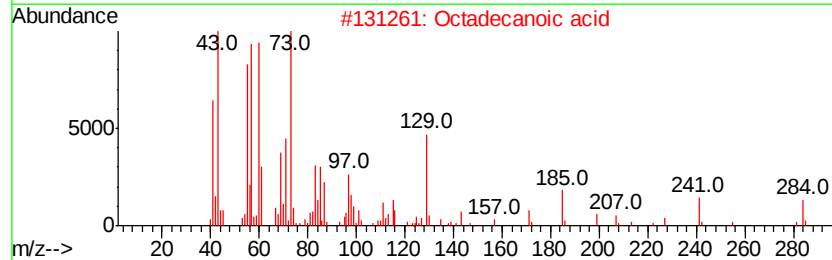
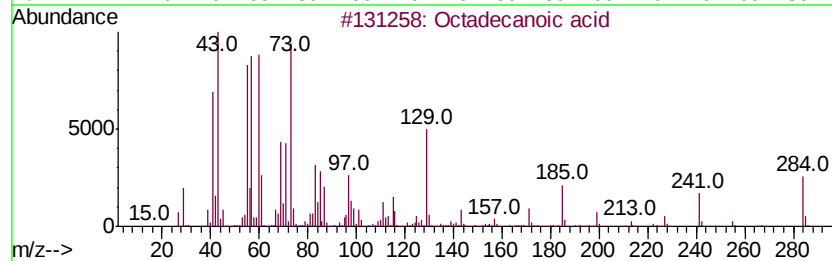
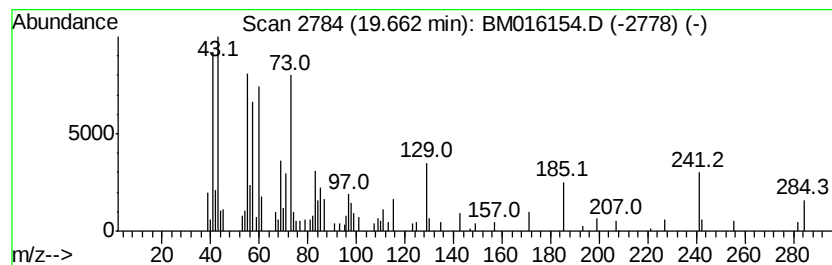
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Peak Number 3 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.66	4.28 ng/ul	245074	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	94
3		Octadecanoic acid	284	C18H36O2	000057-11-4	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	81
5		Octadecanoic acid	284	C18H36O2	000057-11-4	74



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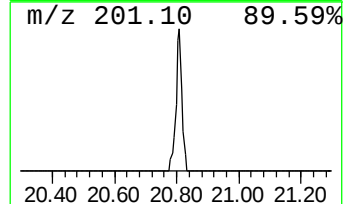
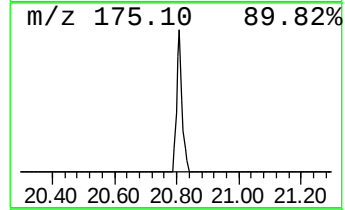
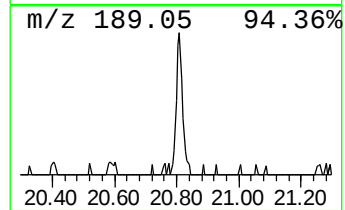
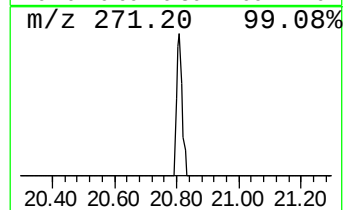
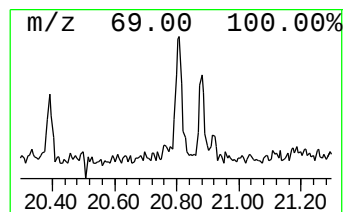
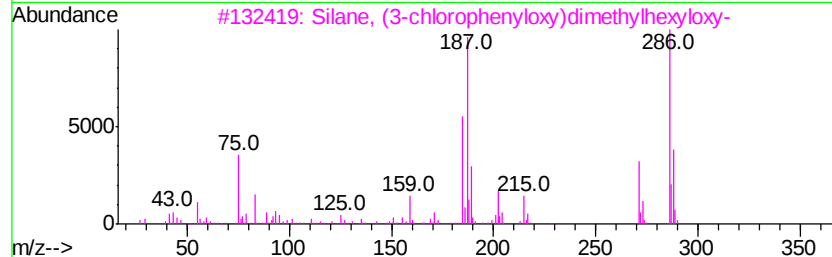
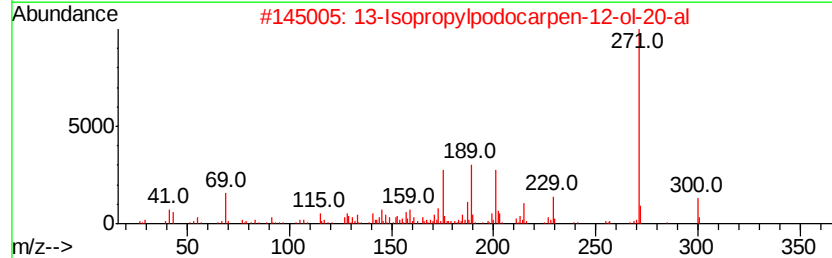
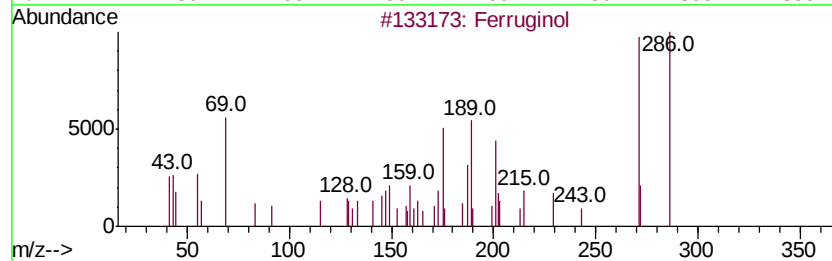
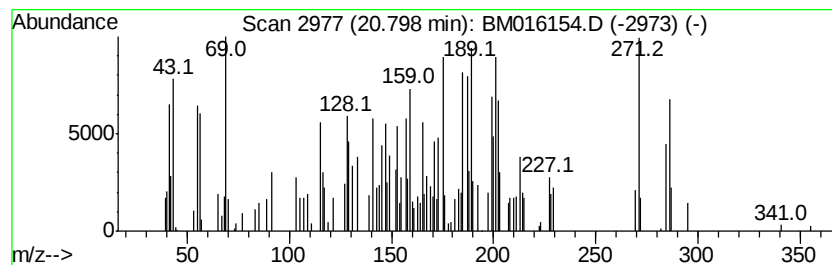
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Peak Number 4 Ferruginol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.80	3.34 ng/ul	190997	Chrysene-d12	21.67
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ferruginol	286	C20H30O	000514-62-5 98
2	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8 47
3	Silane, (3-chlorophenyl)dimethyl-	286	C14H23ClO2Si	1000346-87-9 38
4	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9 30
5	5-Androstene, 4,4-dimethyl-	286	C21H34	1000194-15-4 30



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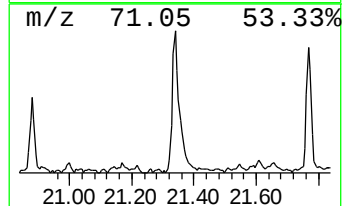
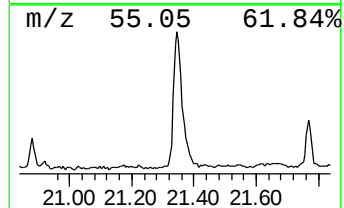
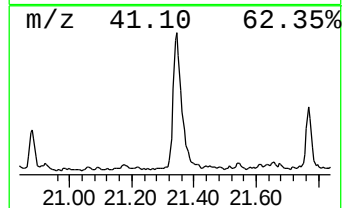
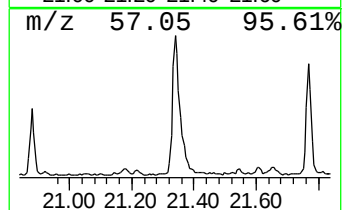
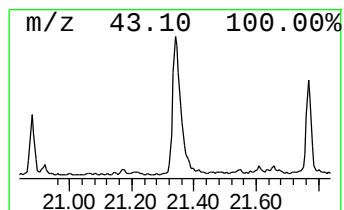
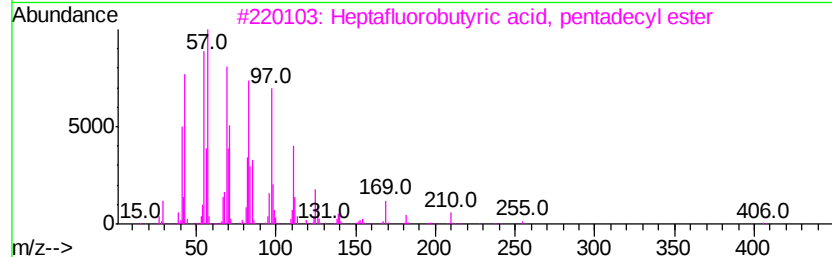
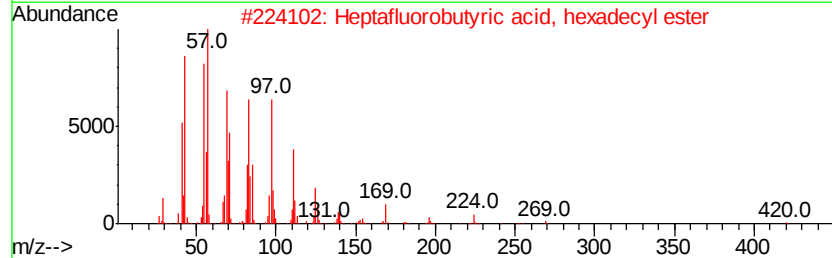
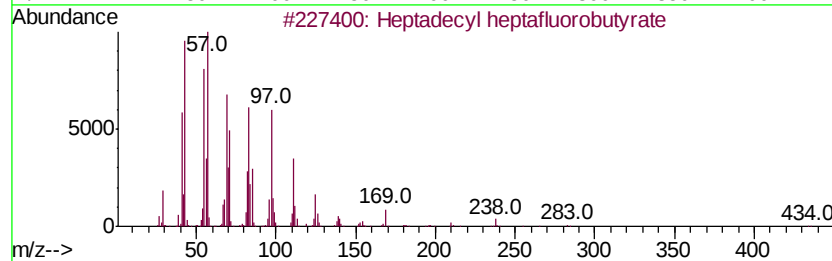
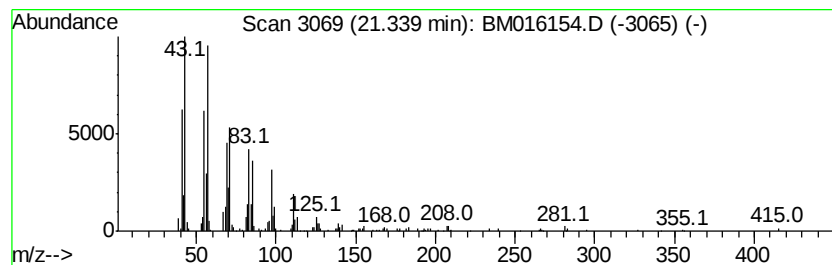
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Heptadecyl heptafluorobutyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.34	9.42 ng/ul	539113	Chrysene-d12	21.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	95
2		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
3		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
4		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	91
5		Docosyl trifluoroacetate	422	C24H45F3O2	1000351-75-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
Operator : SJ/JU
Sample : J4171-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JJ1L4

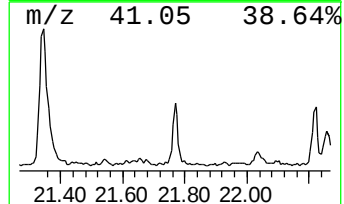
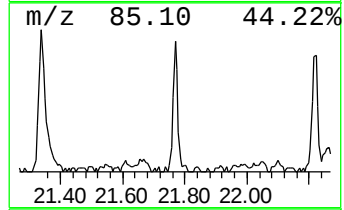
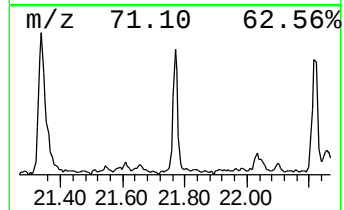
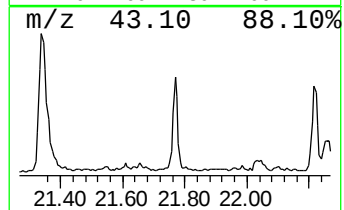
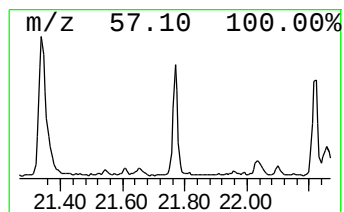
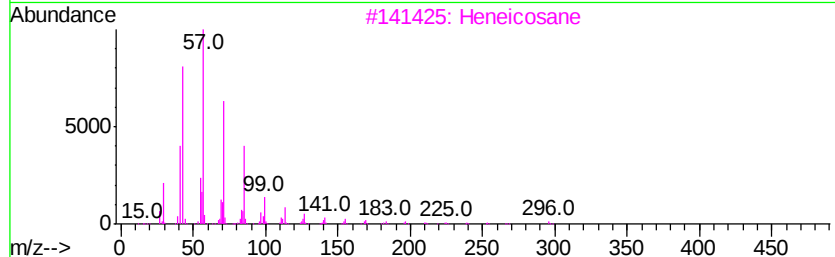
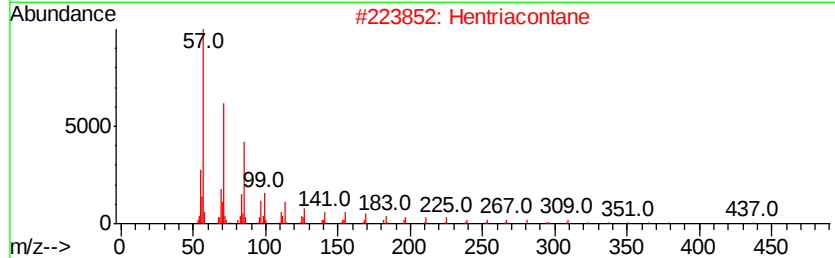
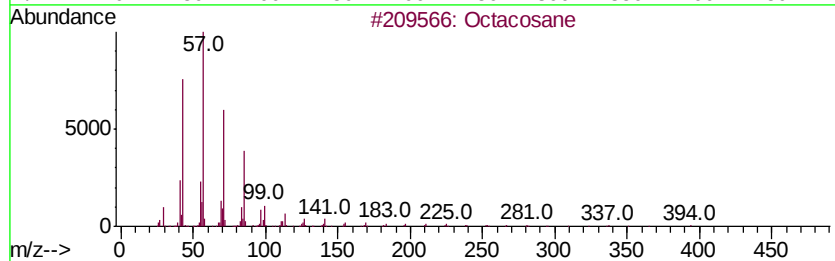
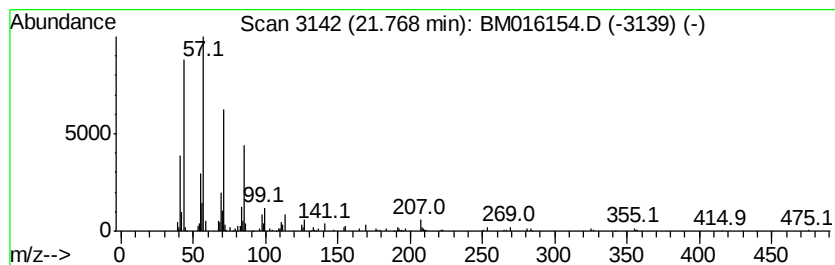
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.77	2.32 ng/ul	132714	Chrysene-d12	21.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	87
2			Hentriacontane	437	C31H64	000630-04-6	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
Operator : SJ/JU
Sample : J4171-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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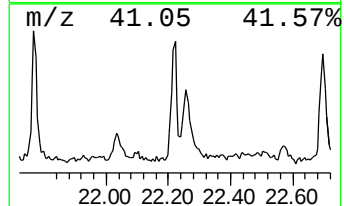
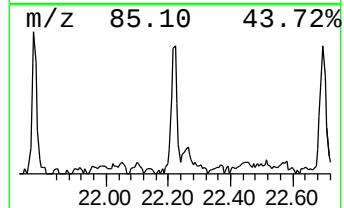
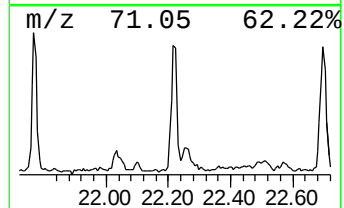
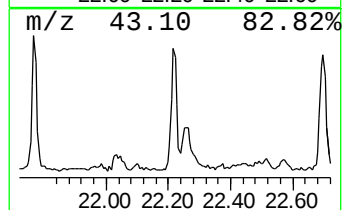
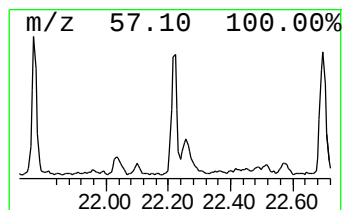
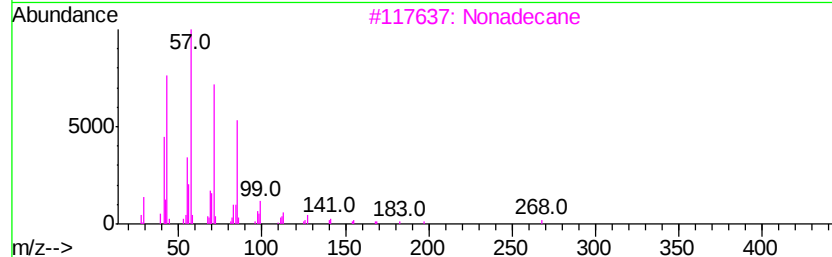
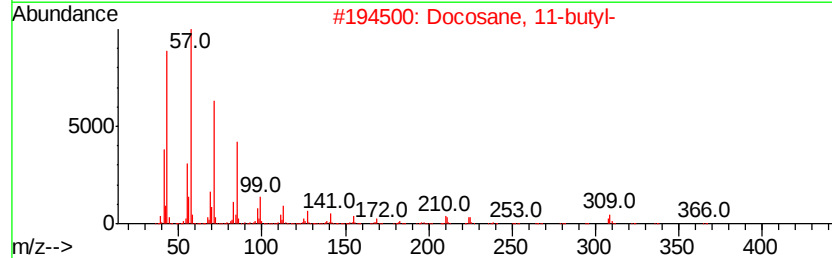
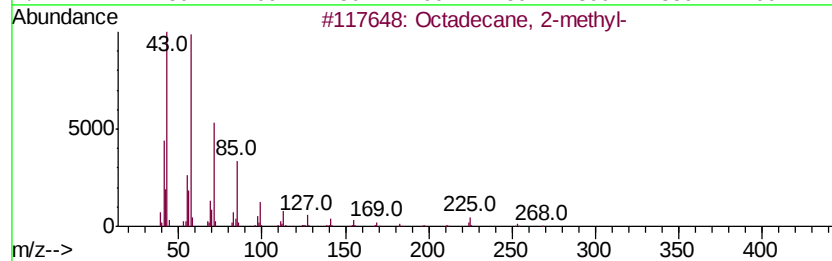
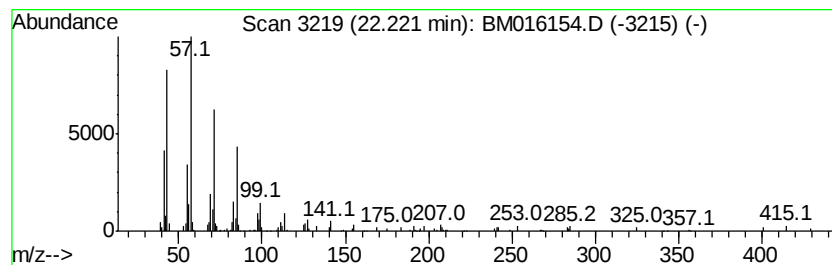
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.22	2.72 ng/ul	155486	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	87
3		Nonadecane	268	C19H40	000629-92-5	87
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
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Sample : J4171-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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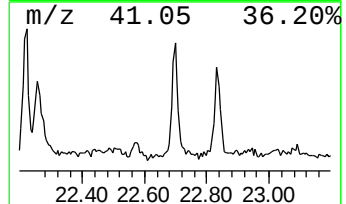
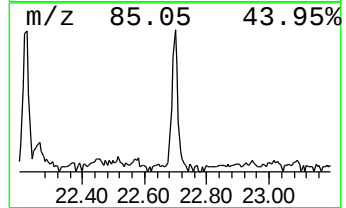
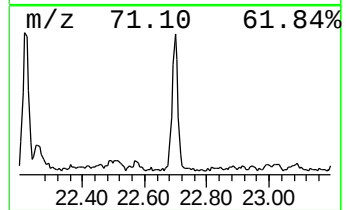
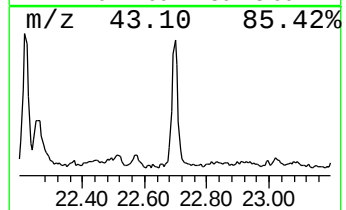
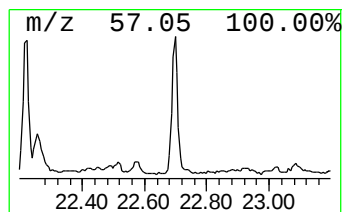
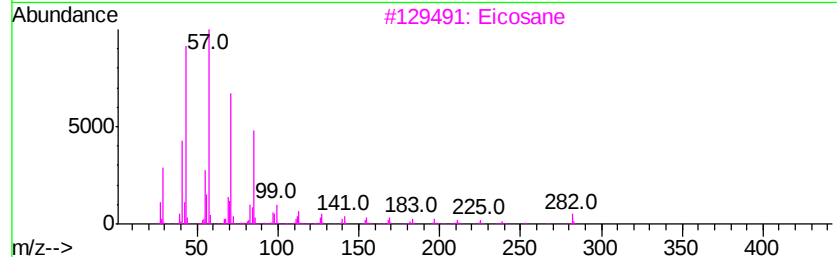
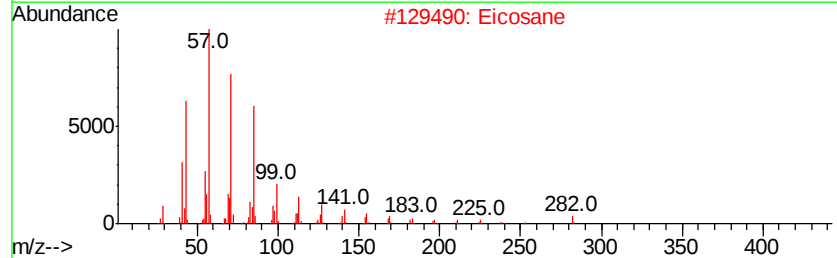
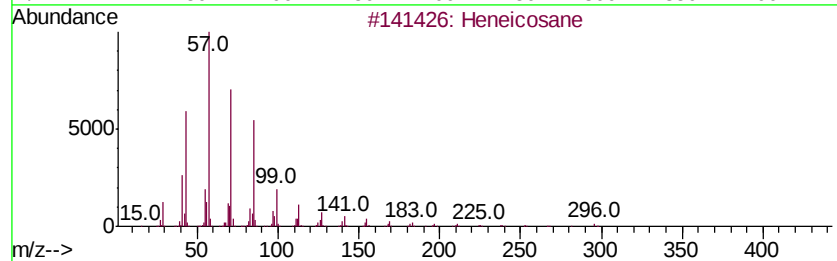
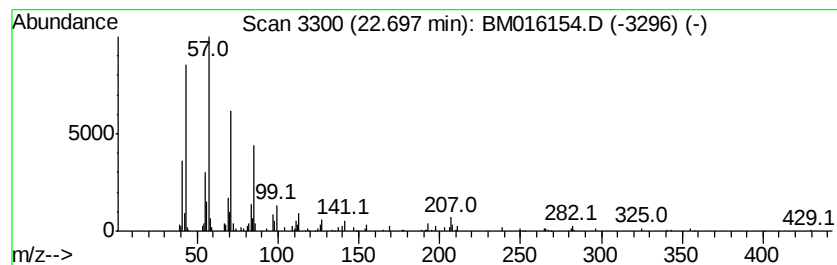
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.70	2.88 ng/ul	164818	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	99
2		Eicosane	282	C20H42	000112-95-8	98
3		Eicosane	282	C20H42	000112-95-8	97
4		Heneicosane	296	C21H44	000629-94-7	97
5		Heneicosane	296	C21H44	000629-94-7	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
Operator : SJ/JU
Sample : J4171-21
Misc :
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Instrument :
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ClientSampleId :
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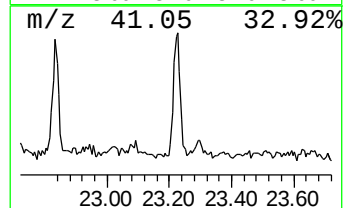
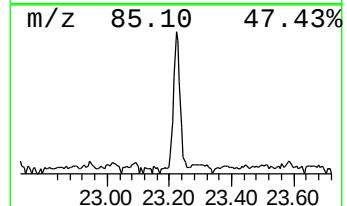
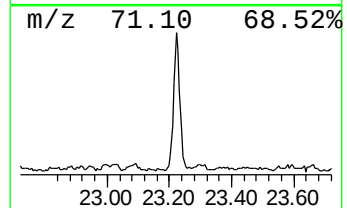
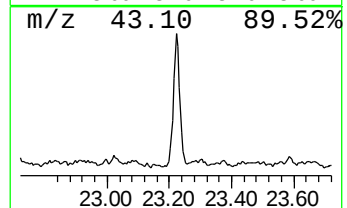
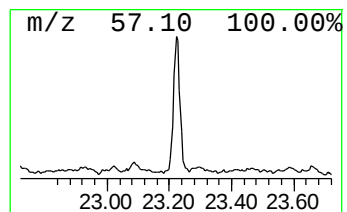
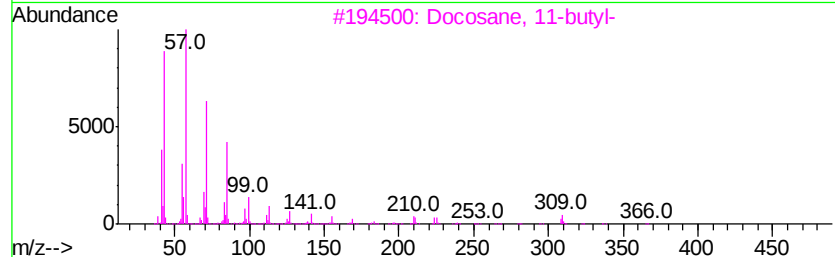
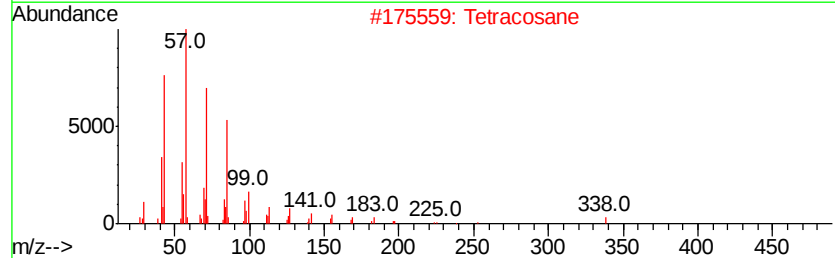
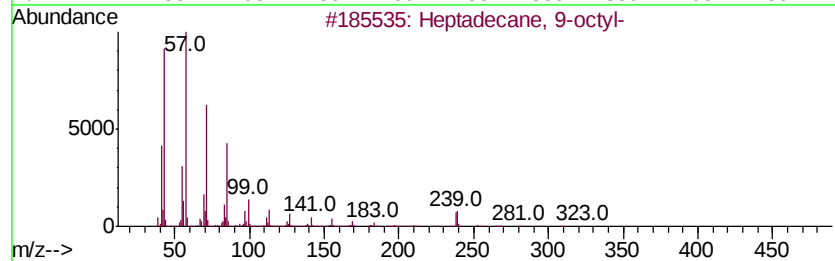
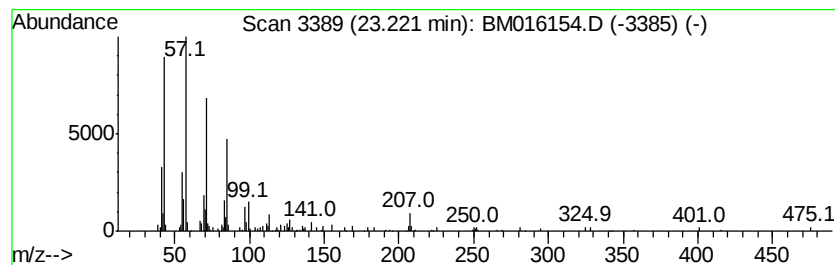
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.22	3.69 ng/ul	177837	Perylene-d12	24.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
2		Tetracosane	338	C24H50	000646-31-1	87
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
Operator : SJ/JU
Sample : J4171-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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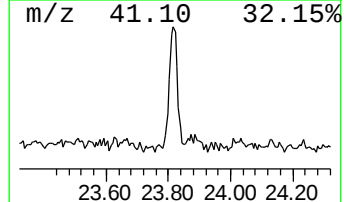
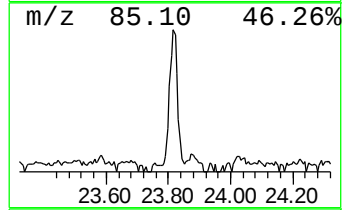
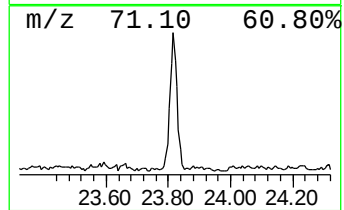
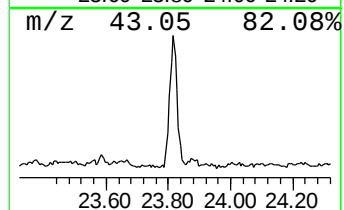
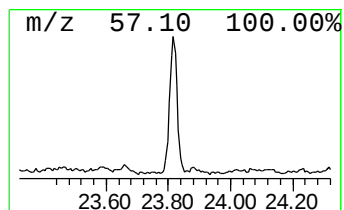
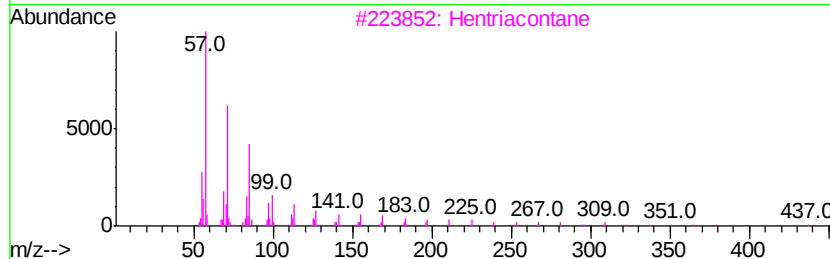
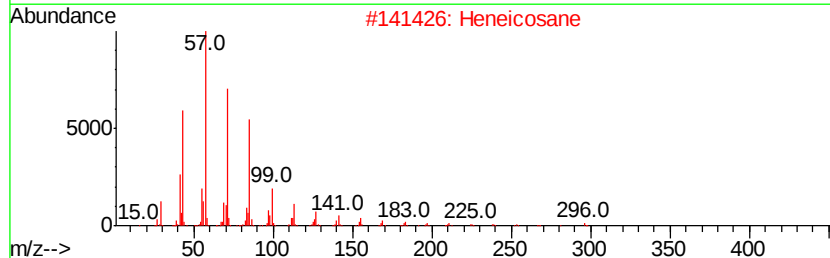
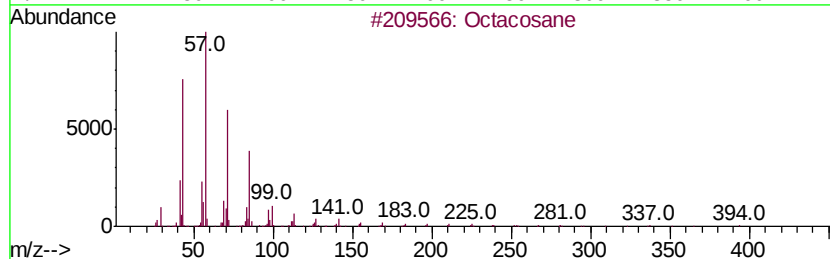
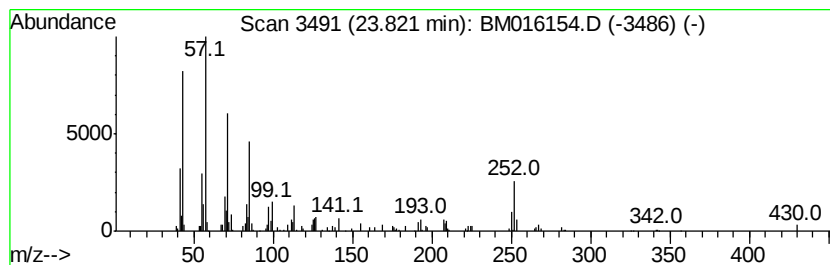
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.82	3.33 ng/ul	160505	Perylene-d12	24.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	87
2			Heneicosane	296	C21H44	000629-94-7	87
3			Hentriacontane	437	C31H64	000630-04-6	87
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080118\
Data File : BM016154.D
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ClientSampleId :
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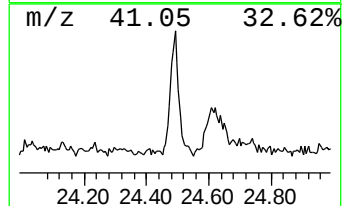
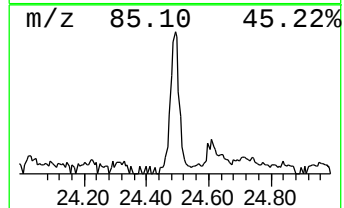
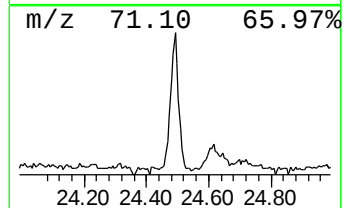
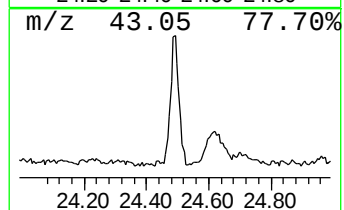
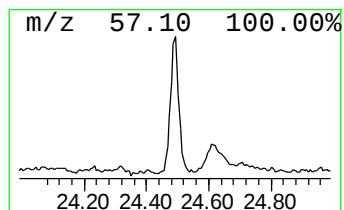
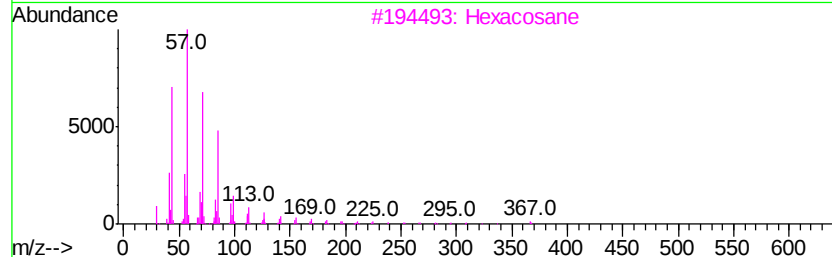
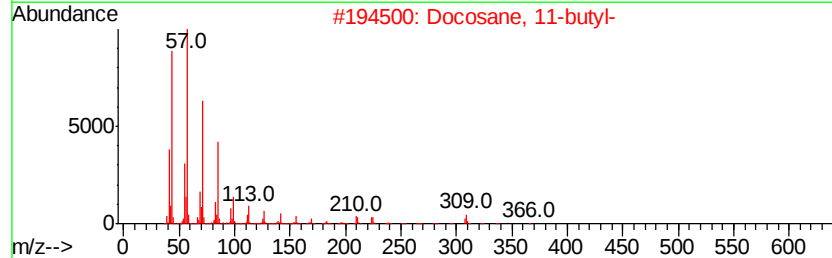
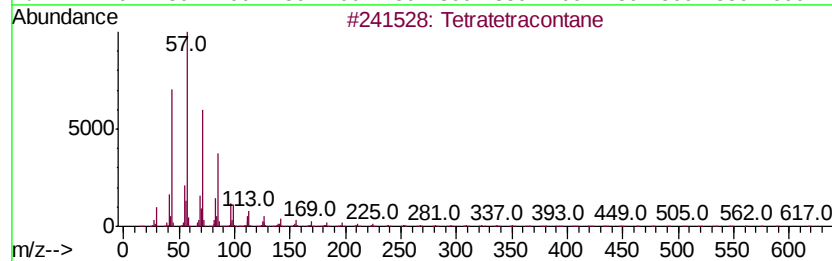
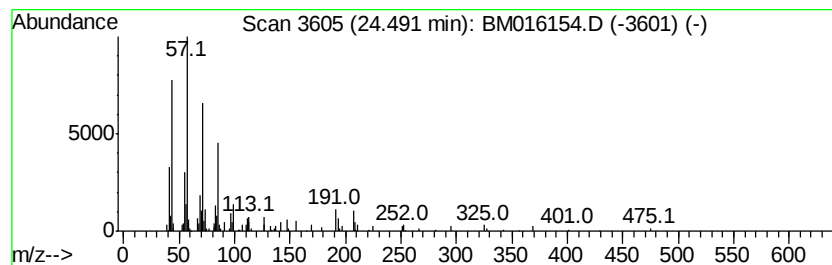
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.49	2.97 ng/ul	142798	Perylene-d12	24.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	76
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	76
3			Hexacosane	366	C26H54	000630-01-3	76
4			Eicosane	282	C20H42	000112-95-8	76
5			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080118\
Data File : BM016154.D
Acq On : 02 Aug 2018 00:16
Operator : SJ/JU
Sample : J4171-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JJ1L4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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n-Hexadecanoic acid	18.40	8.3	ng/ul	382053	4	17.55	921661	20.0
Octadecanoic acid	19.66	4.3	ng/ul	245074	5	21.67	1144020	20.0
Ferruginol	20.80	3.3	ng/ul	190997	5	21.67	1144020	20.0
Heptadecyl heptaf...	21.34	9.4	ng/ul	539113	5	21.67	1144020	20.0
(DEL) Alkane: Str...	21.77	2.3	ng/ul	132714	5	21.67	1144020	20.0
(DEL) Alkane: Str...	22.22	2.7	ng/ul	155486	5	21.67	1144020	20.0
(DEL) Alkane: Str...	22.70	2.9	ng/ul	164818	5	21.67	1144020	20.0
(DEL) Alkane: Str...	23.22	3.7	ng/ul	177837	6	24.03	962801	20.0
(DEL) Alkane: Str...	23.82	3.3	ng/ul	160505	6	24.03	962801	20.0
(DEL) Alkane: Str...	24.49	3.0	ng/ul	142798	6	24.03	962801	20.0