

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
 Data File : BM016946.D
 Acq On : 02 Oct 2018 06:04
 Operator : MJ/SJ
 Sample : J5125-11
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B79

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-BM091018MA.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	3.187	30	34	40	rVB	1305453	1316788	6.76%	1.717%
2	4.787	303	306	314	rVB	83542	113035	0.58%	0.147%
3	7.710	795	803	812	rBV	1362070	2082324	10.69%	2.715%
4	8.928	1002	1010	1015	rBV4	18069	32819	0.17%	0.043%
5	10.269	1233	1238	1244	rVB4	18054	32756	0.17%	0.043%
6	10.351	1244	1252	1259	rVB2	37843	59516	0.31%	0.078%
7	10.487	1262	1275	1284	rBV	1856482	3250312	16.69%	4.237%
8	10.628	1293	1299	1309	rVV7	12780	33685	0.17%	0.044%
9	10.739	1311	1318	1325	rBV6	10033	20862	0.11%	0.027%
10	11.922	1512	1519	1525	rBV	96110	149302	0.77%	0.195%
11	12.228	1565	1571	1577	rBV8	7810	19715	0.10%	0.026%
12	12.416	1598	1603	1607	rVB5	12969	20023	0.10%	0.026%
13	12.781	1659	1665	1669	rVB5	30117	51703	0.27%	0.067%
14	12.833	1669	1674	1679	rBV7	26279	55371	0.28%	0.072%
15	12.886	1679	1683	1687	rVV	64261	87191	0.45%	0.114%
16	12.939	1688	1692	1699	rVB6	25524	47844	0.25%	0.062%
17	13.122	1715	1723	1730	rBV5	73895	187338	0.96%	0.244%
18	13.180	1730	1733	1734	rBV	18616	19987	0.10%	0.026%
19	13.569	1795	1799	1802	rBV5	30796	49065	0.25%	0.064%
20	13.769	1828	1833	1837	rBV	325471	480211	2.47%	0.626%
21	13.839	1841	1845	1849	rVV2	501255	681002	3.50%	0.888%
22	14.004	1869	1873	1877	rBV6	76367	128006	0.66%	0.167%
23	14.051	1878	1881	1883	rBV3	62076	80732	0.41%	0.105%
24	14.092	1883	1888	1891	rBV4	180698	347359	1.78%	0.453%
25	14.180	1899	1903	1911	rVB	580671	991917	5.09%	1.293%
26	14.251	1911	1915	1920	rBV6	196981	387603	1.99%	0.505%
27	14.351	1921	1932	1940	rBV2	2936560	5489730	28.19%	7.157%
28	14.675	1985	1987	1991	rBV	188009	273665	1.41%	0.357%
29	14.763	1999	2002	2006	rBV3	350191	489116	2.51%	0.638%
30	14.886	2018	2023	2028	rBV3	1050404	1669142	8.57%	2.176%
31	15.139	2062	2066	2071	rVB2	1159018	1637374	8.41%	2.135%
32	15.398	2107	2110	2113	rBV4	413194	699059	3.59%	0.911%
33	15.627	2143	2149	2154	rBV	3477660	5590531	28.71%	7.288%
34	15.833	2181	2184	2188	rBV	727732	848733	4.36%	1.106%

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Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	16.110	2224	2231	2236	rBV	12805434	19475014	100.00%	25.389%
36	16.927	2366	2370	2382	rVB	7547345	11966520	61.45%	15.601%
37	17.104	2396	2400	2408	rBV3	2949633	6370850	32.71%	8.306%
38	17.533	2470	2473	2478	rVB	3084638	4321947	22.19%	5.634%
39	21.292	3107	3112	3116	rBV	2795350	3457928	17.76%	4.508%
40	23.521	3485	3491	3499	rVB2	1958389	3689509	18.94%	4.810%

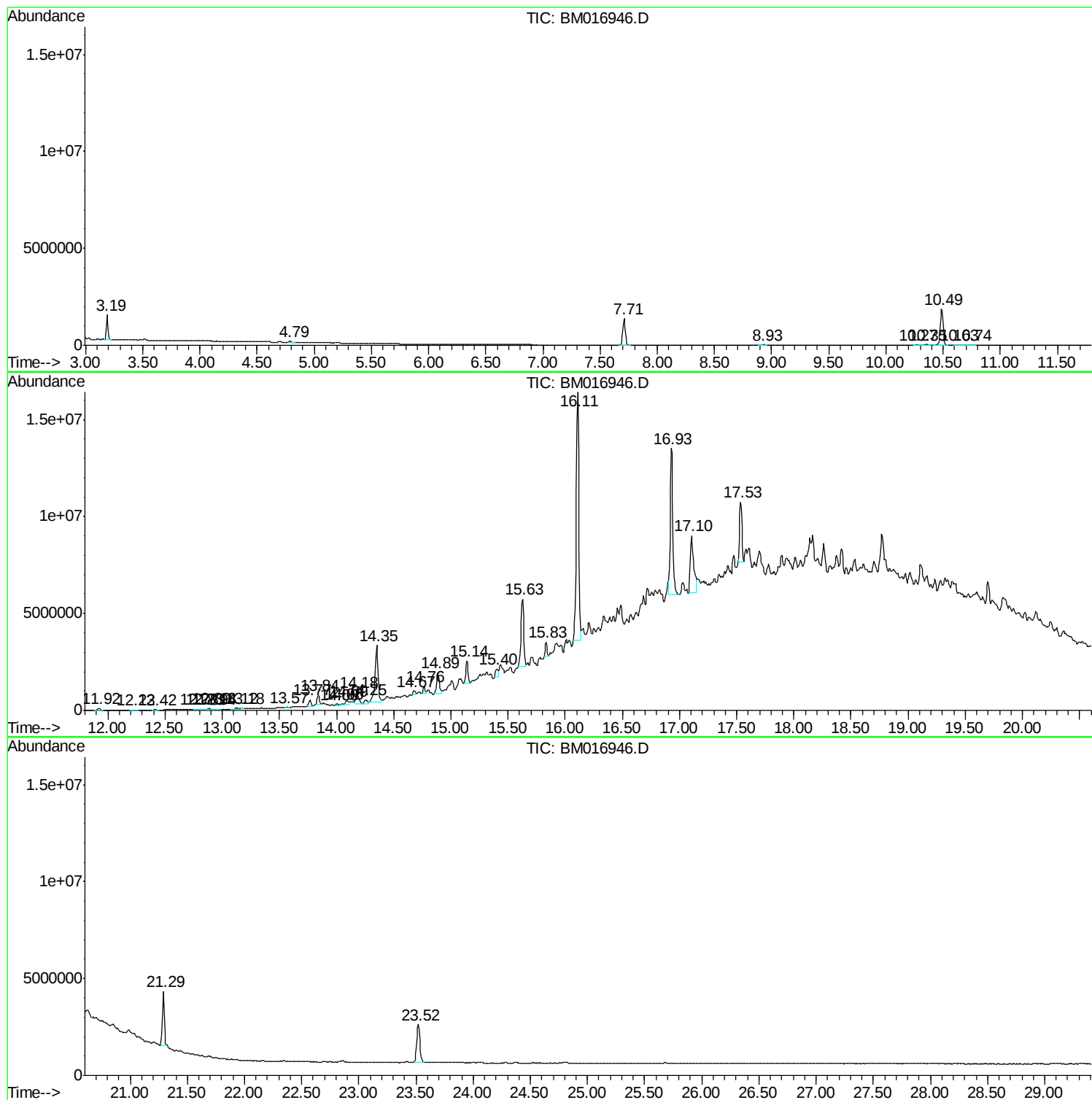
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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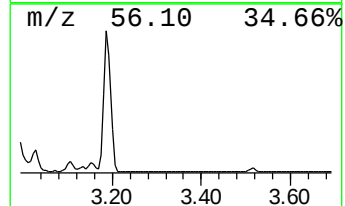
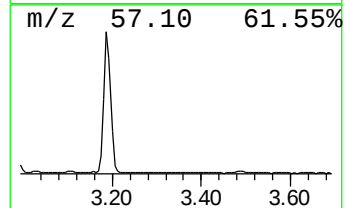
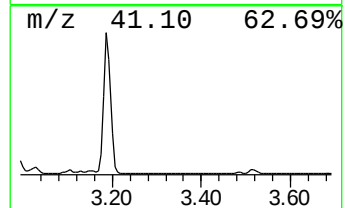
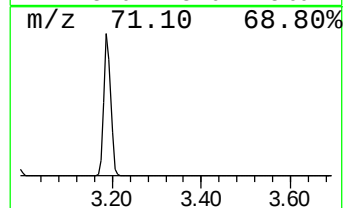
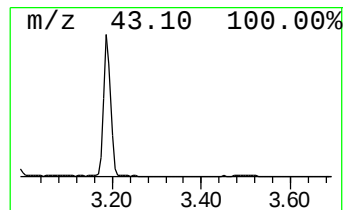
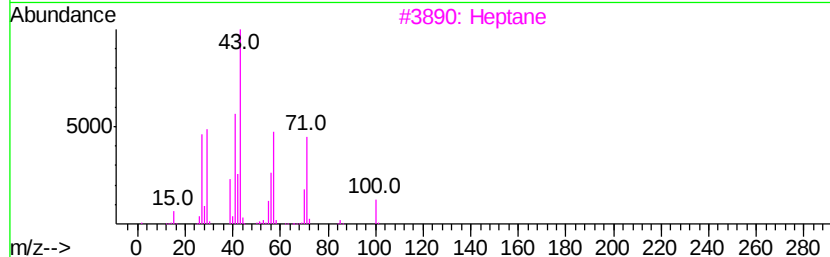
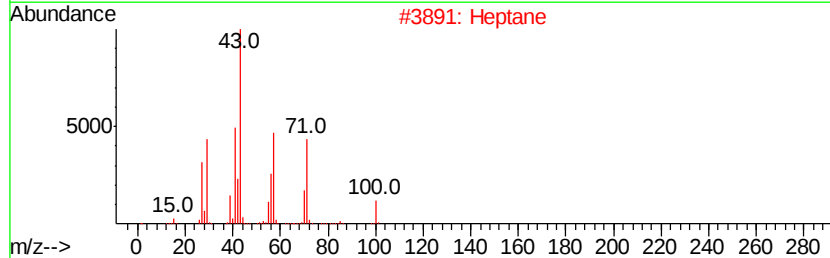
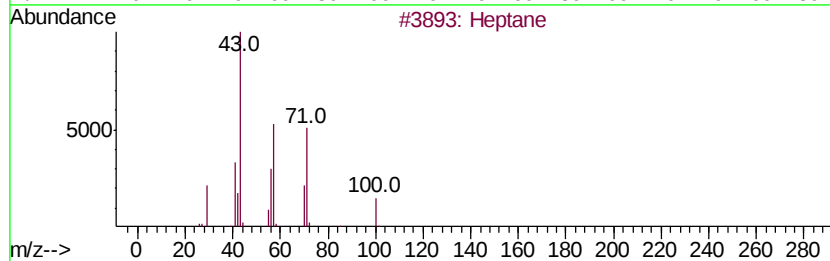
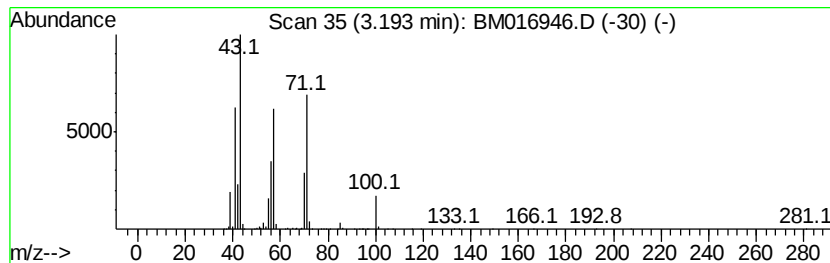
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	12.65 ng/ul	1316790	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	91
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	87
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40



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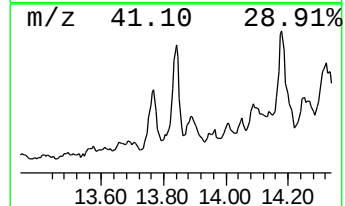
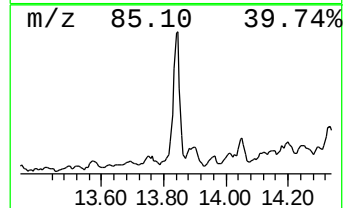
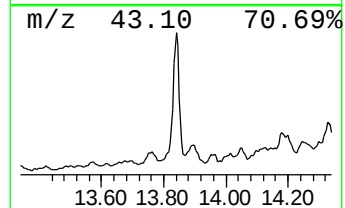
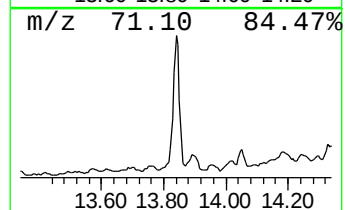
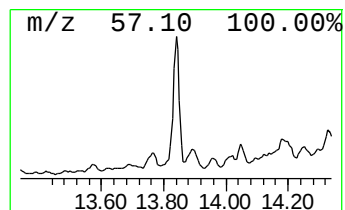
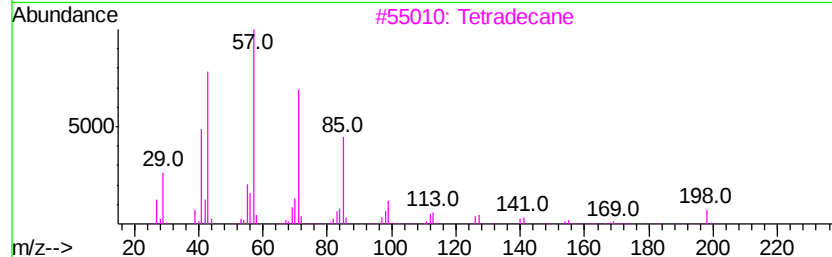
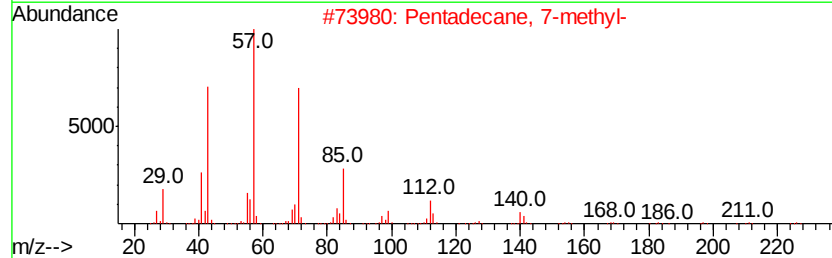
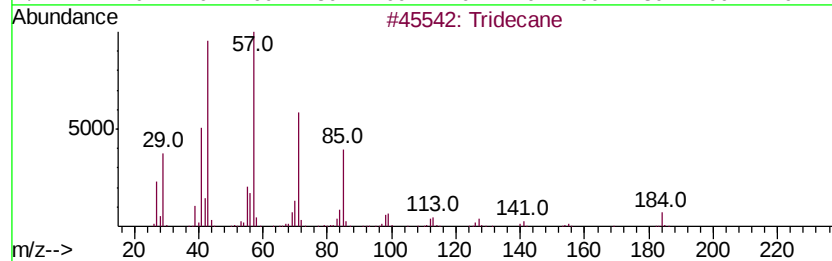
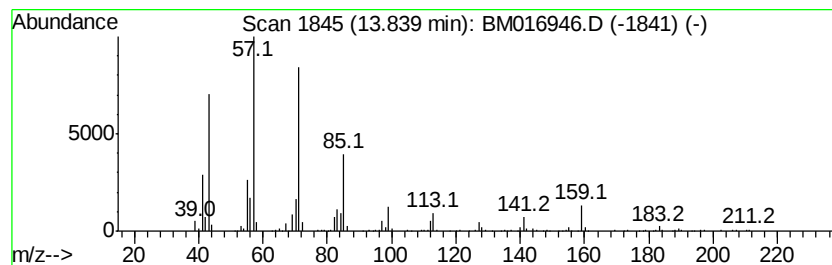
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	2.48 ng/ul	681002	Acenaphthene-d10	14.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	86
2		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	86
3		Tetradecane	198	C14H30	000629-59-4	86
4		Decane, 5-propyl-	184	C13H28	017312-62-8	83
5		Dodecane, 4-methyl-	184	C13H28	006117-97-1	81



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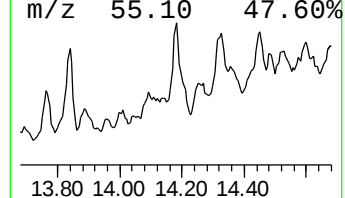
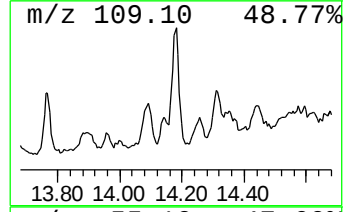
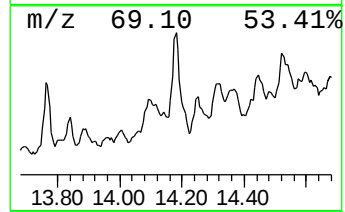
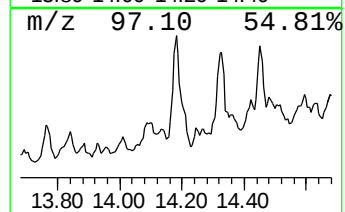
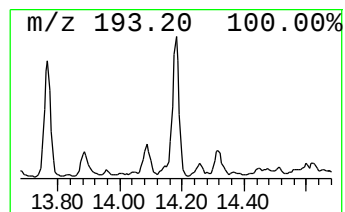
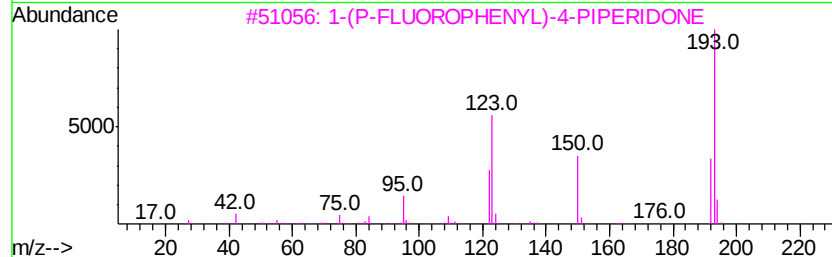
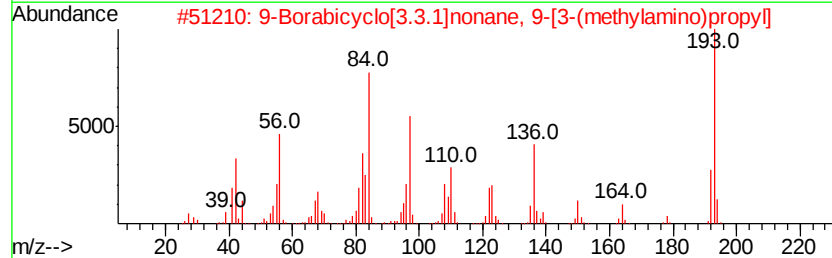
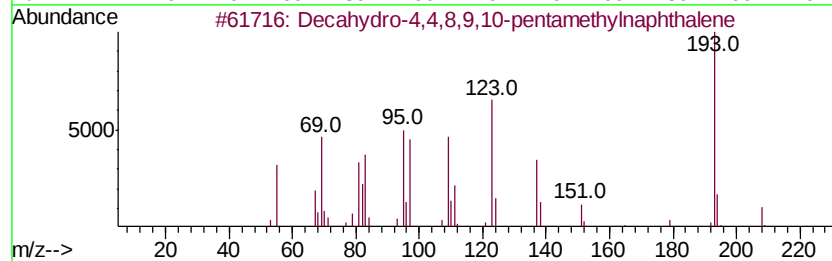
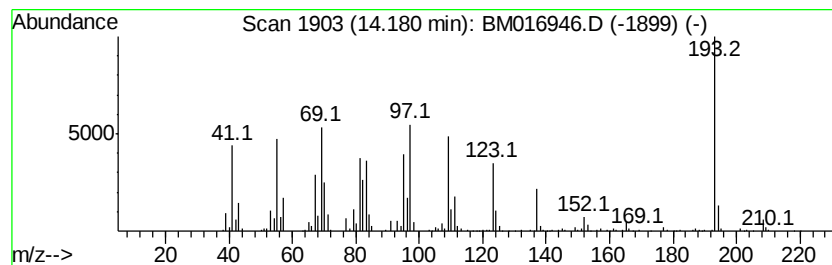
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Decahydro-4,4,8,9,10-pentam... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	3.61 ng/ul	991917	Acenaphthene-d10	14.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3	58
2	9-Borabicyclo[3.3.1]nonane, 9-[3...	193 C12H24BN	1000162-56-0	43
3	1-(P-FLUOROPHENYL)-4-PIPERIDONE	193 C11H12FNO	1000238-56-7	32
4	2-Anthracenamine	193 C14H11N	000613-13-8	27
5	Borinic acid, diethyl-, 3,3,5-tr...	208 C13H25BO	057387-76-5	25



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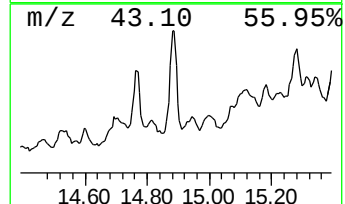
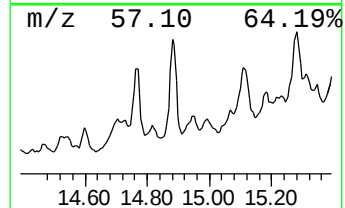
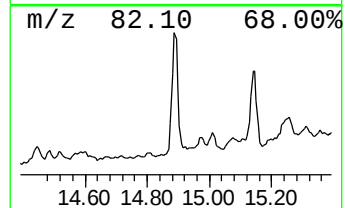
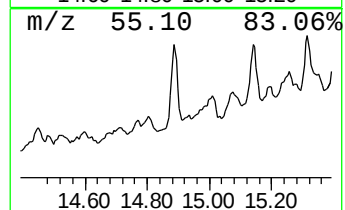
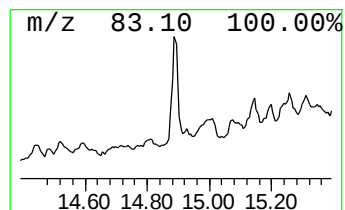
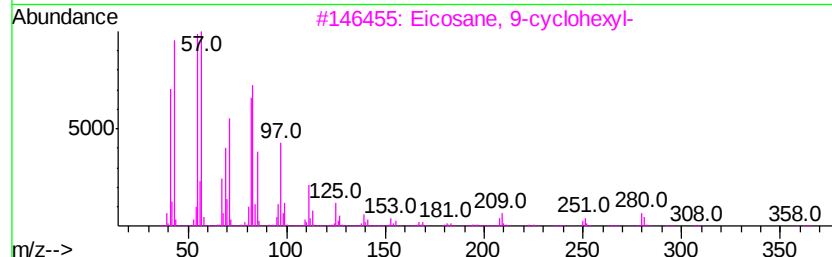
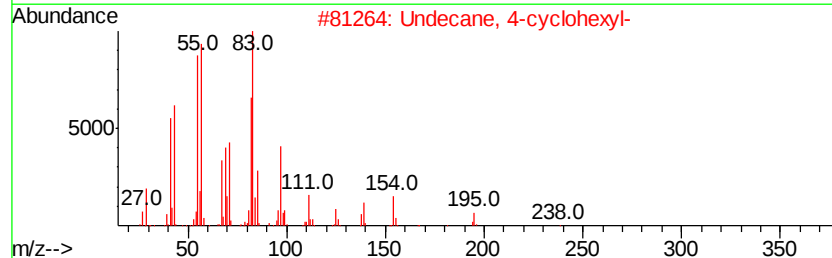
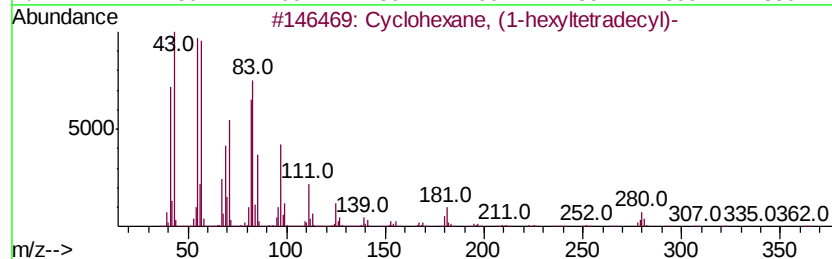
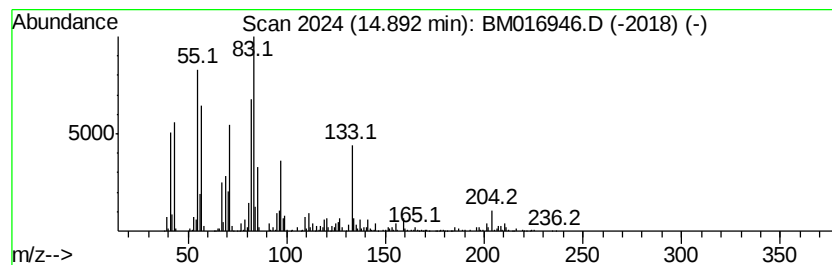
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Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	6.08 ng/ul	1669140	Acenaphthene-d10	14.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, (1-hexyltetradecyl)-	364 C26H52	004443-60-1 49
2		Undecane, 4-cyclohexyl-	238 C17H34	013151-79-6 49
3		Eicosane, 9-cyclohexyl-	364 C26H52	004443-61-2 49
4		Nonadecane	268 C19H40	000629-92-5 25
5		Undecane, 2-methyl-	170 C12H26	007045-71-8 25



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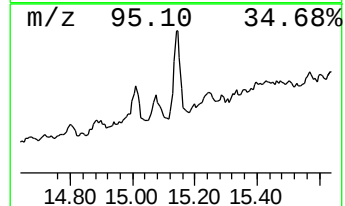
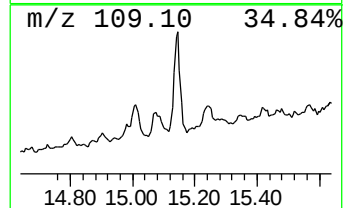
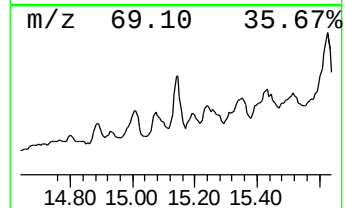
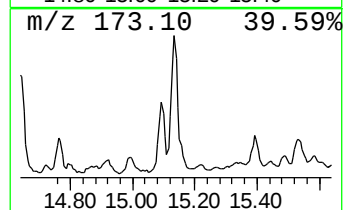
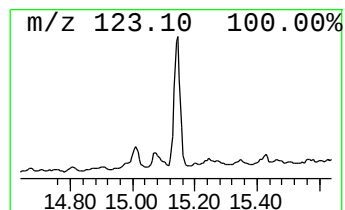
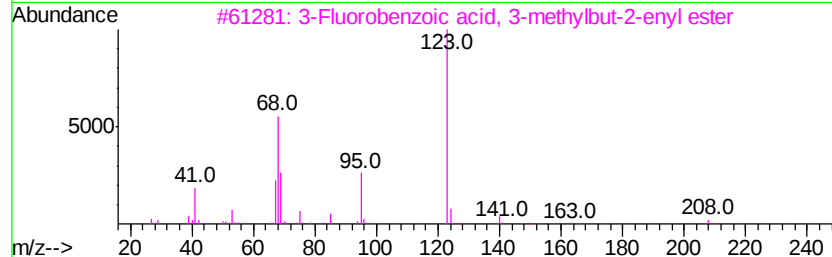
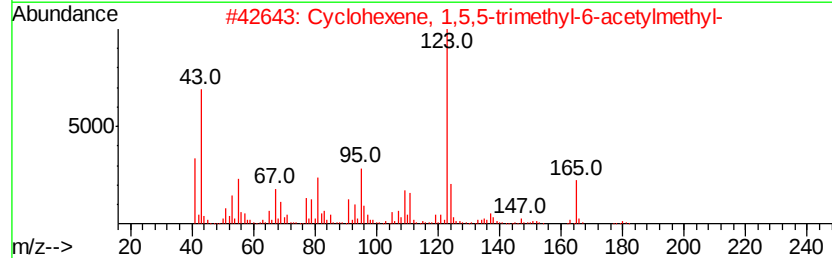
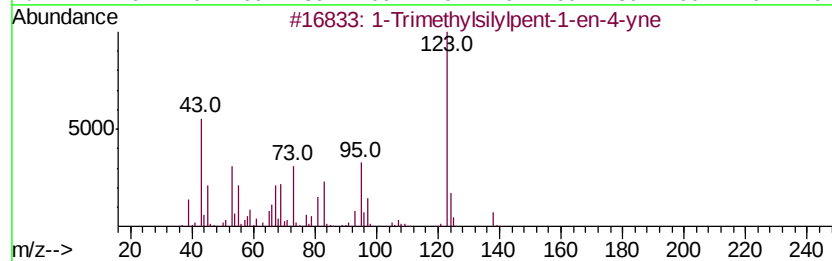
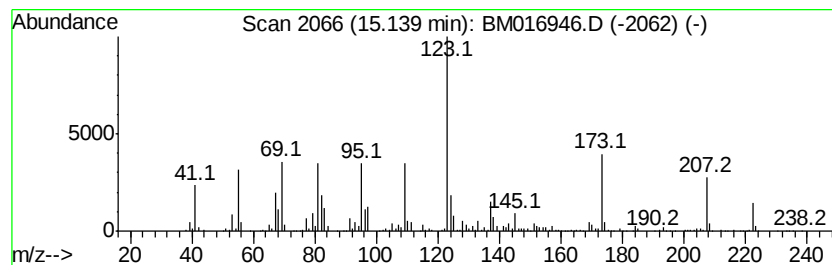
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TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	5.97 ng/ul	1637370	Acenaphthene-d10	14.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	41
2		Cyclohexene, 1,5,5-trimethyl-6-acetyl-	180	C12H20O	211563-96-1	38
3		3-Fluorobenzoic acid, 3-methylbut-2-enyl ester	208	C12H13FO2	154559-38-3	35
4		4-Chloro-2-isopropyl-5-methylphenol	306	C18H23ClO2	321918-37-0	35
5		Bicyclo[2.2.2]octane, 1-methyl-4-ylidene-	202	C10H18O2S	069855-48-7	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
Operator : MJ/SJ
Sample : J5125-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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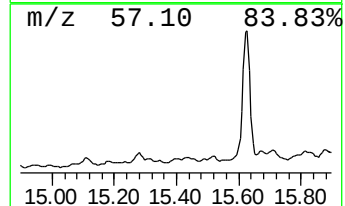
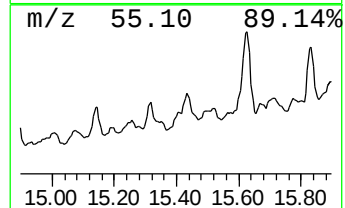
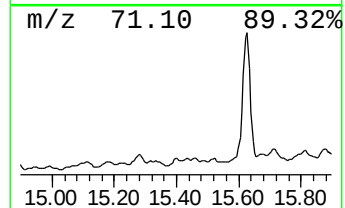
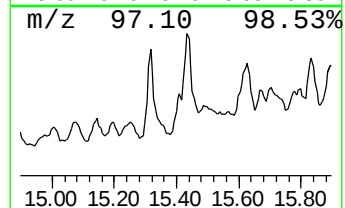
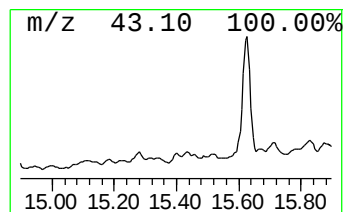
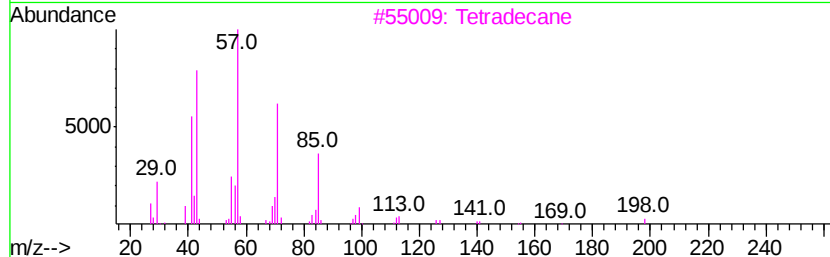
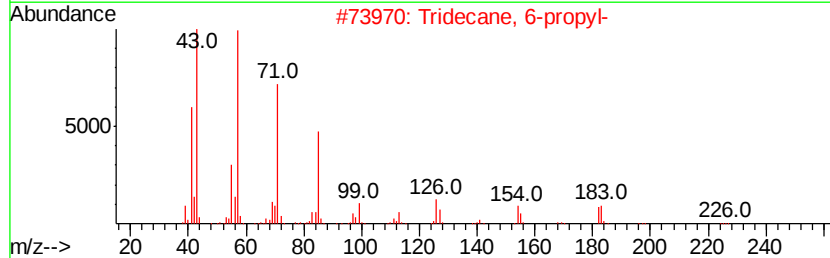
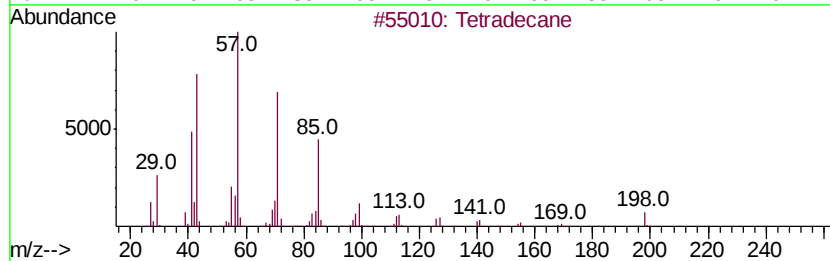
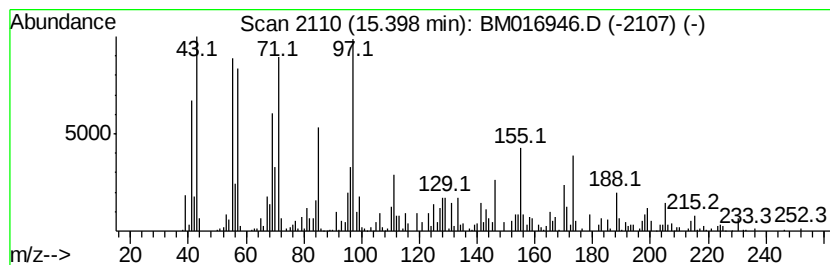
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	2.55 ng/ul	699059	Acenaphthene-d10	14.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	47
2			Tridecane, 6-propyl-	226	C16H34	055045-10-8	38
3			Tetradecane	198	C14H30	000629-59-4	35
4			Tetradecane	198	C14H30	000629-59-4	25
5			Diisoamylene	140	C10H20	054063-09-1	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
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Sample : J5125-11
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Instrument :
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ClientSampleId :
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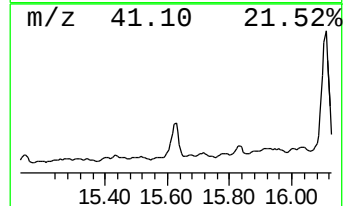
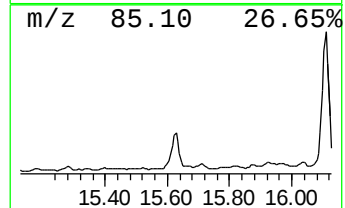
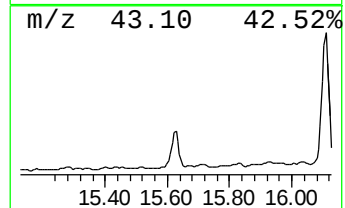
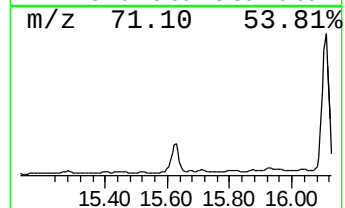
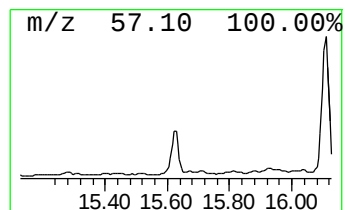
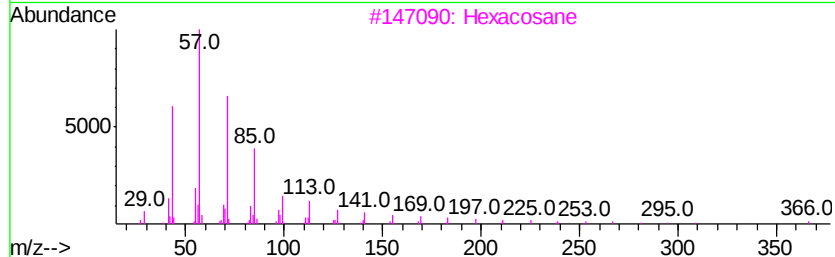
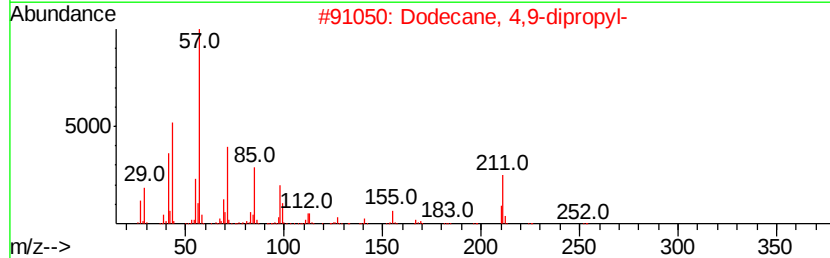
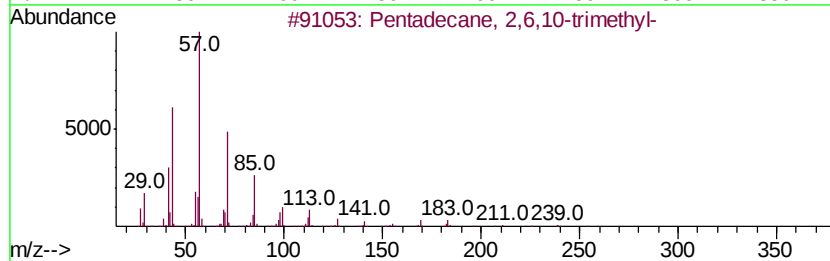
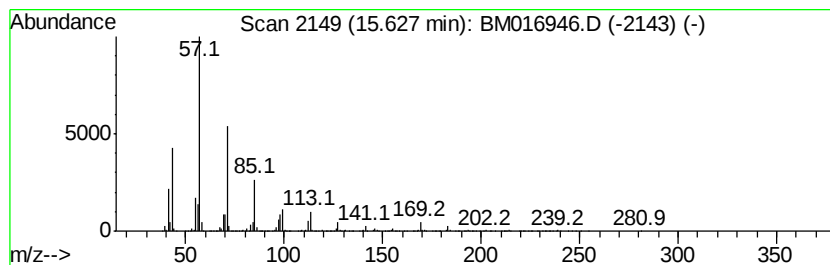
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.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.
15.63	20.37 ng/ul	5590530	Acenaphthene-d10	14.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	94
2		Dodecane, 4,9-dipropyl-	254	C18H38	003054-63-5	90
3		Hexacosane	366	C26H54	000630-01-3	86
4		Hexadecane	226	C16H34	000544-76-3	80
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
Operator : MJ/SJ
Sample : J5125-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

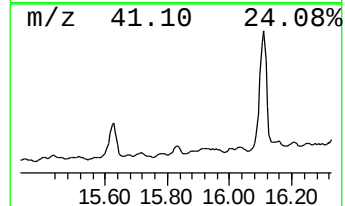
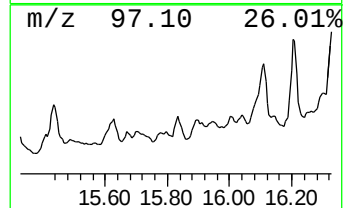
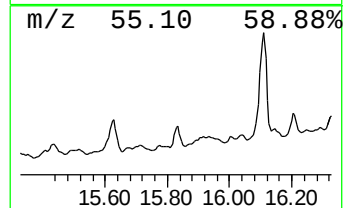
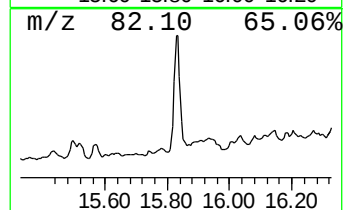
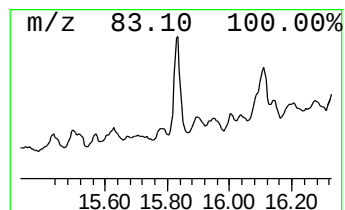
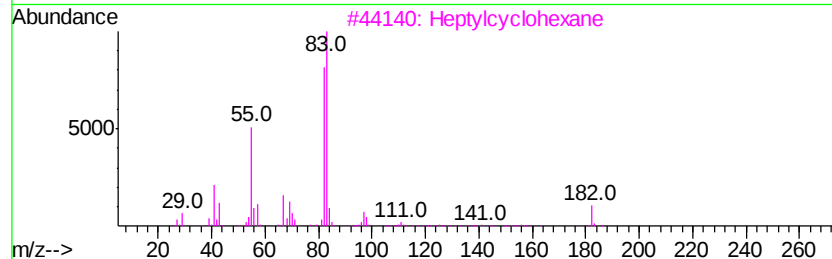
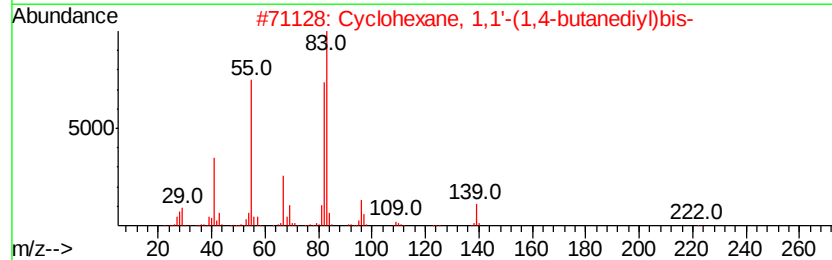
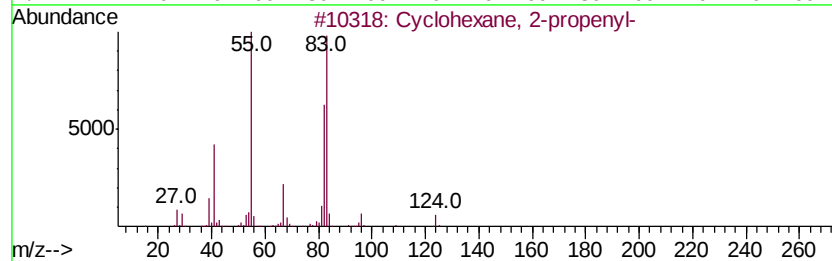
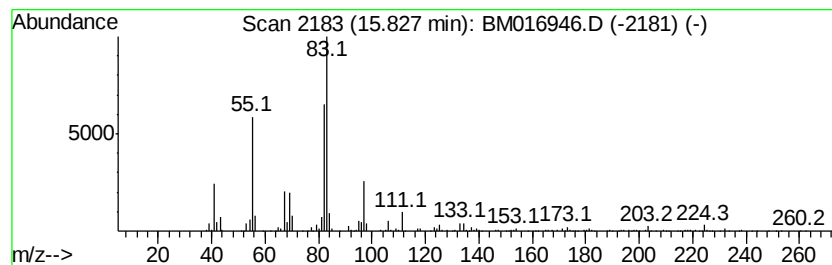
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Cyclohexane, 2-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	2.66 ng/ul	848733	Phenanthrene-d10	17.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 2-propenyl-	124 C9H16	002114-42-3 70
2		Cyclohexane, 1,1'-(1,4-butanediyl)bis-	222 C16H30	006165-44-2 70
3		Heptylcyclohexane	182 C13H26	005617-41-4 58
4		Cyclohexane, decyl-	224 C16H32	001795-16-0 58
5		n-Amylcyclohexane	154 C11H22	029949-27-7 53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
Operator : MJ/SJ
Sample : J5125-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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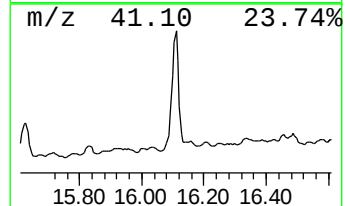
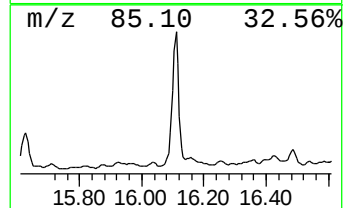
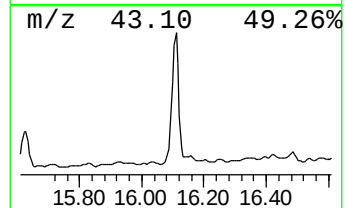
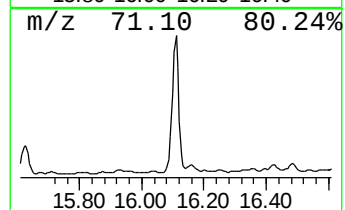
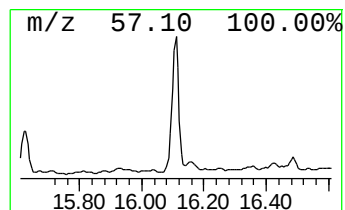
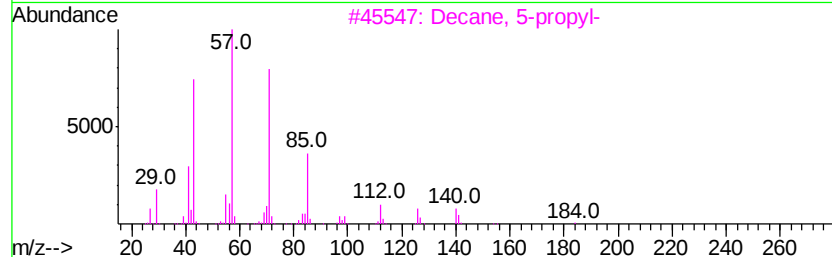
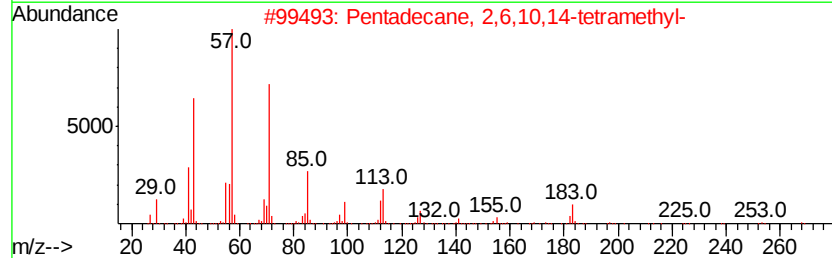
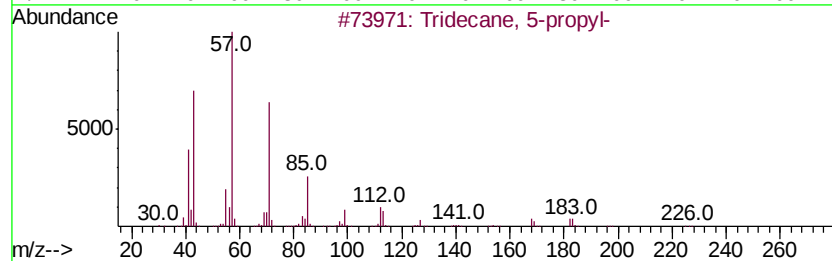
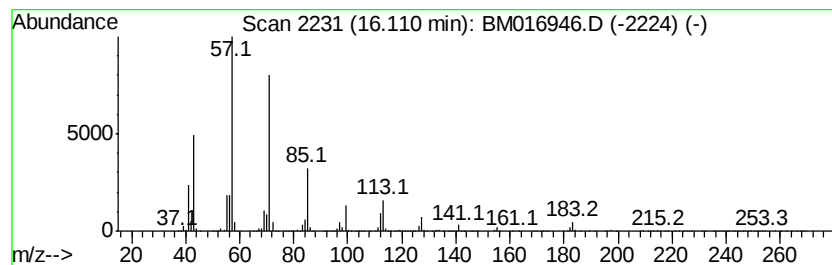
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.11	61.14 ng/ul	19475000	Phenanthrene-d10	17.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
3		Decane, 5-propyl-	184	C13H28	017312-62-8	83
4		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
5		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
Operator : MJ/SJ
Sample : J5125-11
Misc :
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Instrument :
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ClientSampleId :
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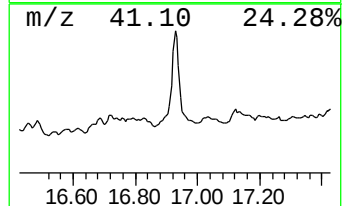
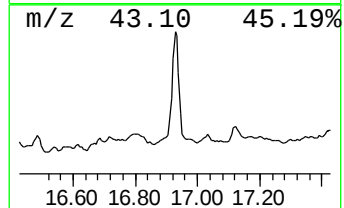
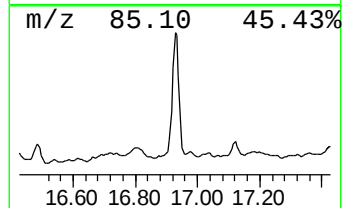
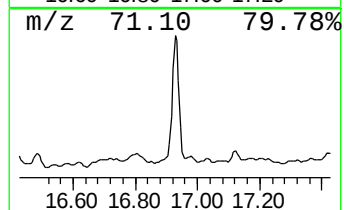
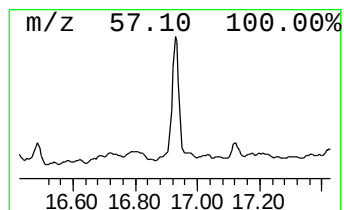
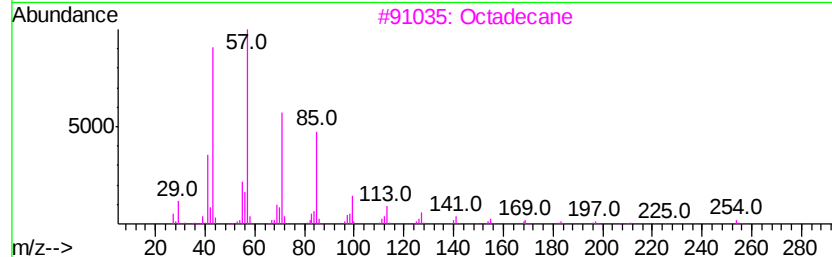
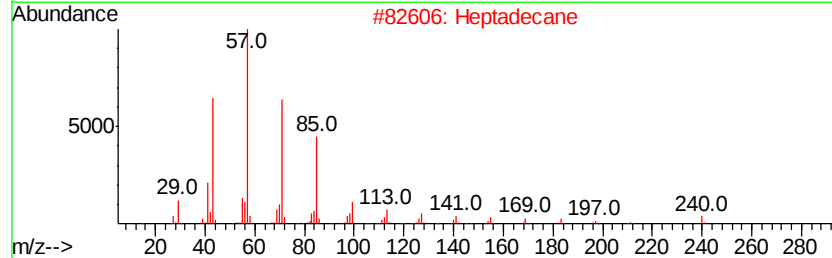
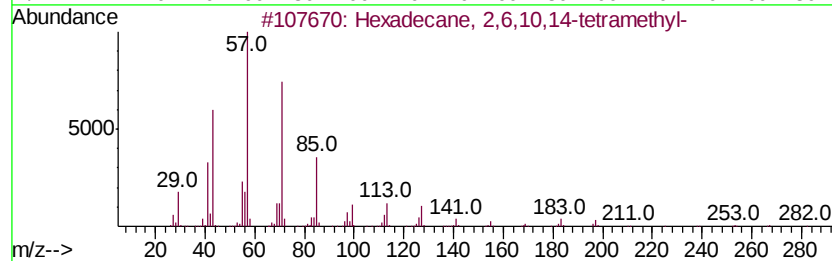
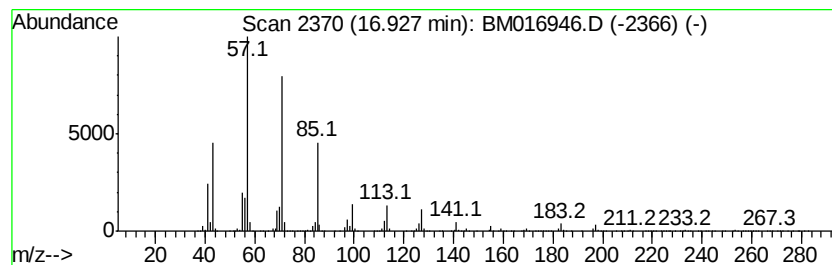
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.93	37.57 ng/ul	11966500	Phenanthrene-d10	17.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Octadecane	254	C18H38	000593-45-3	90
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
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Acq On : 02 Oct 2018 06:04
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Sample : J5125-11
Misc :
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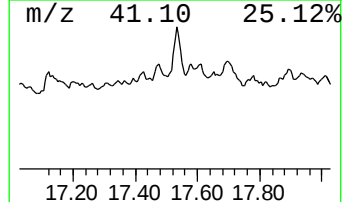
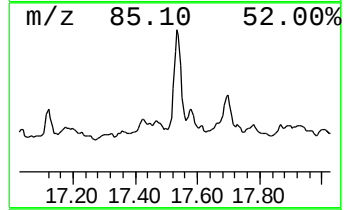
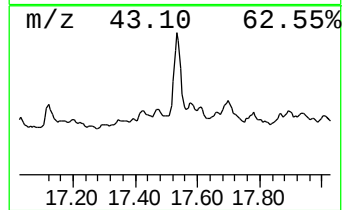
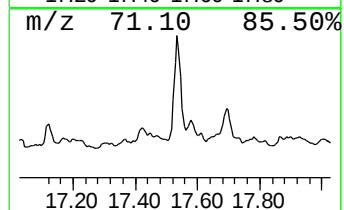
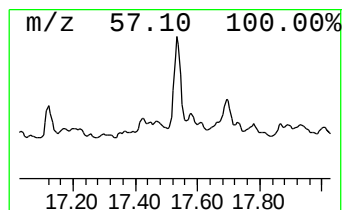
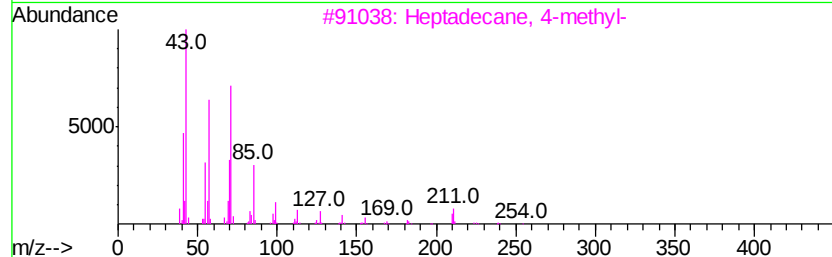
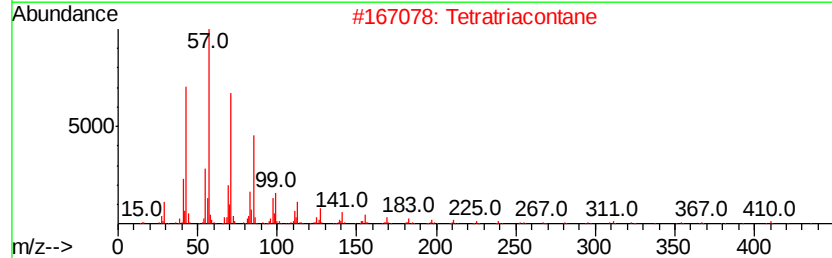
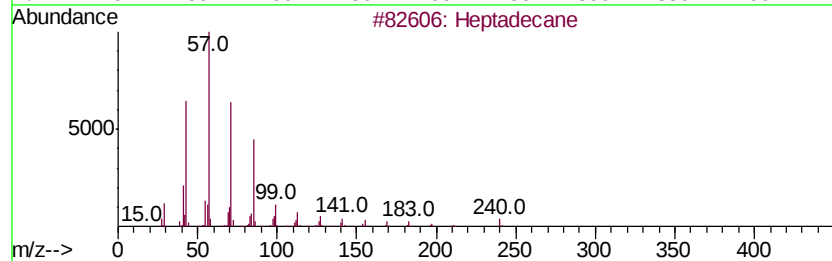
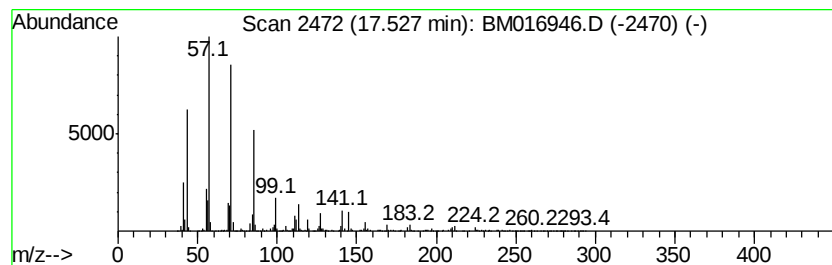
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.53	13.57 ng/ul	4321950	Phenanthrene-d10	17.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Tetratriacontane	479	C34H70	014167-59-0	91
3		Heptadecane, 4-methyl-	254	C18H38	026429-11-8	90
4		Heneicosane	296	C21H44	000629-94-7	87
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100118\
Data File : BM016946.D
Acq On : 02 Oct 2018 06:04
Operator : MJ/SJ
Sample : J5125-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: Str...	13.84	2.5	ng/ul	681002	3	14.35	5489730	20.0
Decahydro-4,4,8,9...	14.18	3.6	ng/ul	991917	3	14.35	5489730	20.0
unknown-01	14.89	6.1	ng/ul	1669140	3	14.35	5489730	20.0
unknown-02	15.14	6.0	ng/ul	1637370	3	14.35	5489730	20.0
(DEL) Alkane: Str...	15.40	2.5	ng/ul	699059	3	14.35	5489730	20.0
(DEL) Alkane: Str...	15.63	20.4	ng/ul	5590530	3	14.35	5489730	20.0
Cyclohexane, 2-pr...	15.83	2.7	ng/ul	848733	4	17.10	6370850	20.0
(DEL) Alkane: Str...	16.11	61.1	ng/ul	19475000	4	17.10	6370850	20.0
(DEL) Alkane: Str...	16.93	37.6	ng/ul	11966500	4	17.10	6370850	20.0
(DEL) Alkane: Str...	17.53	13.6	ng/ul	4321950	4	17.10	6370850	20.0