

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
 Data File : BM014700.D
 Acq On : 13 Apr 2018 13:43
 Operator : SJ/JU
 Sample : J2313-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1A39

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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032818MA.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	3.693	13	18	25	rVB	128016	179135	1.38%	0.117%
2	7.004	572	581	592	rBV	138968	252284	1.94%	0.165%
3	7.687	692	697	710	rVB2	500208	1102862	8.50%	0.720%
4	7.869	721	728	737	rBV	1770692	2819336	21.73%	1.839%
5	8.087	758	765	773	rBV	1883986	3059281	23.58%	1.996%
6	8.569	837	847	864	rBV	1850348	3198810	24.66%	2.087%
7	8.951	907	912	920	rVB	107023	179343	1.38%	0.117%
8	9.063	923	931	936	rBV2	89484	154174	1.19%	0.101%
9	9.263	959	965	976	rVB	1477920	2414350	18.61%	1.575%
10	9.692	1027	1038	1042	rBV4	169441	378220	2.92%	0.247%
11	9.751	1042	1048	1063	rVB	2246732	3954821	30.49%	2.580%
12	10.051	1093	1099	1105	rVB2	179646	325963	2.51%	0.213%
13	10.122	1105	1111	1118	rBV2	173796	329600	2.54%	0.215%
14	10.216	1123	1127	1135	rVB	94525	151253	1.17%	0.099%
15	10.480	1165	1172	1185	rBV	1684559	2995376	23.09%	1.954%
16	10.586	1185	1190	1195	rVV3	74337	136103	1.05%	0.089%
17	10.651	1195	1201	1214	rVB2	330540	646702	4.99%	0.422%
18	10.798	1216	1226	1235	rBV2	884257	1580356	12.18%	1.031%
19	11.022	1257	1264	1271	rVV	2588505	4398914	33.91%	2.870%
20	11.422	1323	1332	1344	rVB	2611199	4966385	38.29%	3.240%
21	11.533	1344	1351	1357	rBV3	242907	531525	4.10%	0.347%
22	11.598	1357	1362	1369	rVB3	107569	247769	1.91%	0.162%
23	11.710	1376	1381	1386	rBV	125943	204707	1.58%	0.134%
24	11.763	1386	1390	1397	rVB2	137839	239599	1.85%	0.156%
25	12.145	1449	1455	1461	rBV	178523	312153	2.41%	0.204%
26	12.310	1477	1483	1492	rVB3	138573	277987	2.14%	0.181%
27	12.498	1509	1515	1524	rBV3	138773	337749	2.60%	0.220%
28	12.598	1524	1532	1537	rBV5	60591	138306	1.07%	0.090%
29	12.698	1544	1549	1556	rVB	102773	159000	1.23%	0.104%
30	12.939	1583	1590	1596	rBV3	215621	472624	3.64%	0.308%
31	13.157	1622	1627	1631	rBV	90565	141915	1.09%	0.093%
32	13.251	1632	1643	1651	rVB	8018651	12971561	100.00%	8.463%
33	13.427	1667	1673	1676	rBV4	87286	155936	1.20%	0.102%
34	13.474	1677	1681	1687	rVV4	47986	136194	1.05%	0.089%

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
 Data File : BM014700.D
 Acq On : 13 Apr 2018 13:43
 Operator : SJ/JU
 Sample : J2313-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1A39

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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032818MA.M
 Title : SVOA CALIBRATION

35	13.651	1707	1711	1715	rBV2	114062	196473	1.51%	0.128%
36	13.727	1720	1724	1729	rVV2	90323	168524	1.30%	0.110%
37	13.780	1730	1733	1741	rVB4	72308	143021	1.10%	0.093%
38	14.145	1790	1795	1800	rVB3	134630	241240	1.86%	0.157%
39	14.204	1800	1805	1809	rBV2	176707	269903	2.08%	0.176%
40	14.492	1847	1854	1859	rBV4	265626	619824	4.78%	0.404%
41	14.557	1859	1865	1870	rVV	4478818	6244117	48.14%	4.074%
42	14.604	1870	1873	1877	rVV	432367	616653	4.75%	0.402%
43	14.645	1877	1880	1887	rVB3	110999	193851	1.49%	0.126%
44	14.868	1912	1918	1926	rBV	5345166	8126313	62.65%	5.302%
45	15.021	1940	1944	1949	rVB	214076	289908	2.23%	0.189%
46	15.168	1963	1969	1975	rBV2	3446035	5429145	41.85%	3.542%
47	15.233	1975	1980	1986	rVB	1689015	2521474	19.44%	1.645%
48	15.315	1986	1994	1999	rBV3	424953	863746	6.66%	0.564%
49	15.510	2025	2027	2035	rVB4	100800	185918	1.43%	0.121%
50	15.598	2039	2042	2046	rBV4	99951	173247	1.34%	0.113%
51	15.692	2054	2058	2063	rVB5	135212	252897	1.95%	0.165%
52	15.757	2065	2069	2075	rVB7	137511	258630	1.99%	0.169%
53	15.868	2084	2088	2093	rBV	320498	504661	3.89%	0.329%
54	15.962	2100	2104	2107	rBV	229449	272738	2.10%	0.178%
55	16.145	2129	2135	2143	rBV	6309628	9154682	70.58%	5.973%
56	16.239	2146	2151	2160	rVV	1743050	2576985	19.87%	1.681%
57	16.474	2188	2191	2203	rBV6	196212	461051	3.55%	0.301%
58	16.809	2245	2248	2252	rVB	182580	194020	1.50%	0.127%
59	17.039	2284	2287	2291	rBV2	129523	157731	1.22%	0.103%
60	17.662	2389	2393	2395	rBV5	157586	272973	2.10%	0.178%
61	17.880	2425	2430	2439	rVB2	3881034	5611002	43.26%	3.661%
62	17.980	2441	2447	2456	rVB2	5978347	8638652	66.60%	5.636%
63	18.715	2567	2572	2579	rBV	2255567	3246959	25.03%	2.118%
64	19.839	2759	2763	2770	rBV2	597544	898508	6.93%	0.586%
65	19.950	2778	2782	2802	rBV	3128846	4373018	33.71%	2.853%
66	20.168	2817	2819	2821	rBV	173738	152830	1.18%	0.100%
67	20.215	2821	2827	2837	rVB	5696668	8079356	62.29%	5.271%
68	20.686	2904	2907	2913	rVB	397798	452689	3.49%	0.295%
69	21.168	2986	2989	2996	rVB	739754	866389	6.68%	0.565%
70	21.621	3062	3066	3075	rVB	1484716	1744300	13.45%	1.138%
71	21.962	3119	3124	3131	rVB	3238178	4517305	34.82%	2.947%

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

Integration Parameters: LSCINT.P

Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032818MA.M
Title : SVOA CALIBRATION

72	22.074	3139	3143	3148	rVB	1388349	1625099	12.53%	1.060%
73	22.556	3220	3225	3227	rBV	1278045	1893289	14.60%	1.235%
74	22.586	3227	3230	3239	rVB	2895153	3867430	29.81%	2.523%
75	23.074	3308	3313	3328	rVB	1263227	1978556	15.25%	1.291%
76	23.644	3406	3410	3416	rVB	976533	1532643	11.82%	1.000%
77	24.315	3516	3524	3536	rVB3	3654538	8433122	65.01%	5.502%
78	24.480	3545	3552	3562	rVB	2051321	4312948	33.25%	2.814%
79	25.050	3644	3649	3660	rVB	530165	1116068	8.60%	0.728%

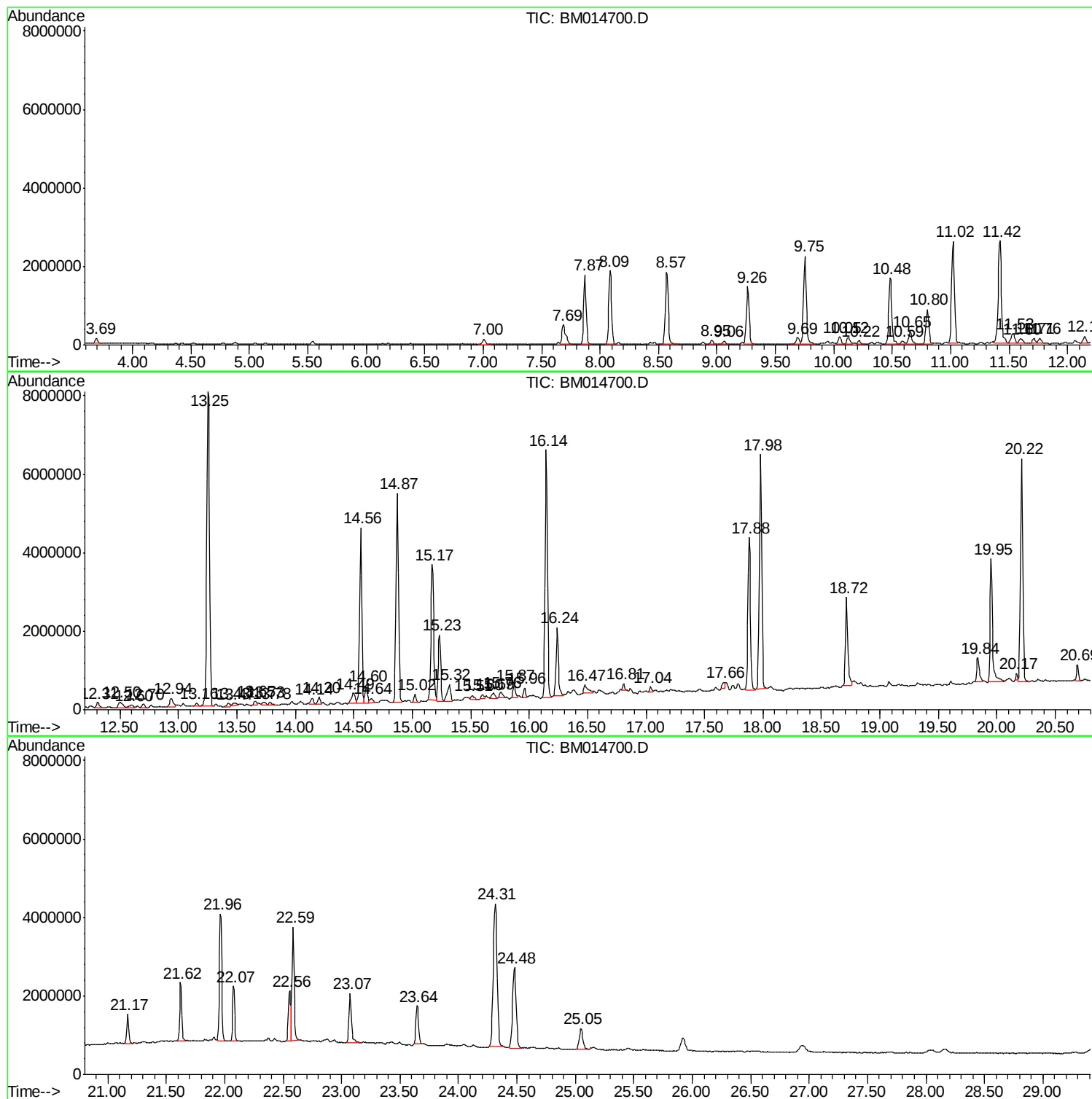
Sum of corrected areas: 153280181

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

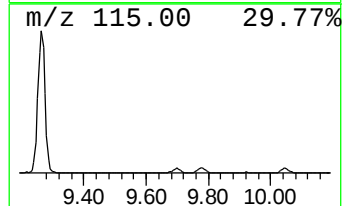
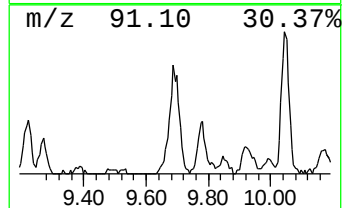
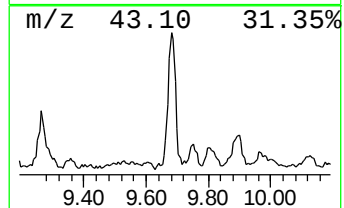
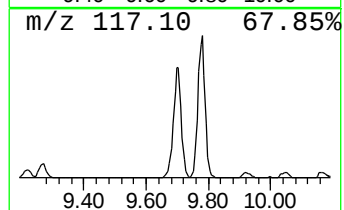
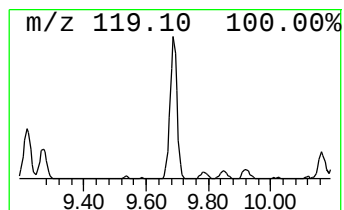
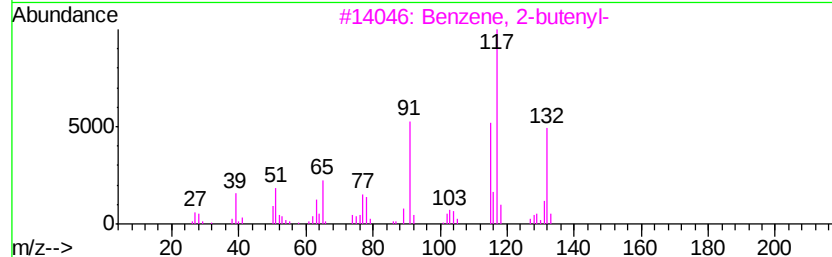
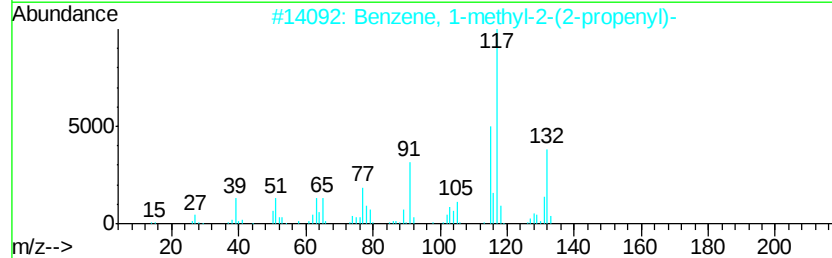
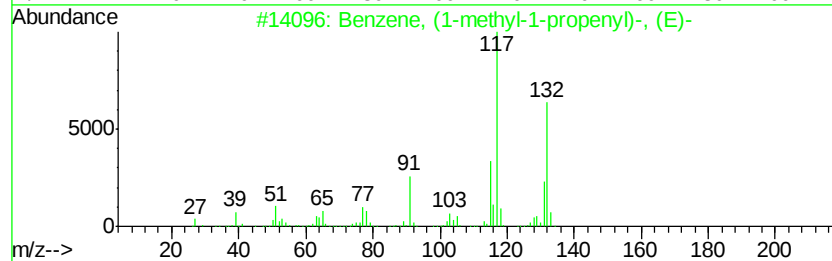
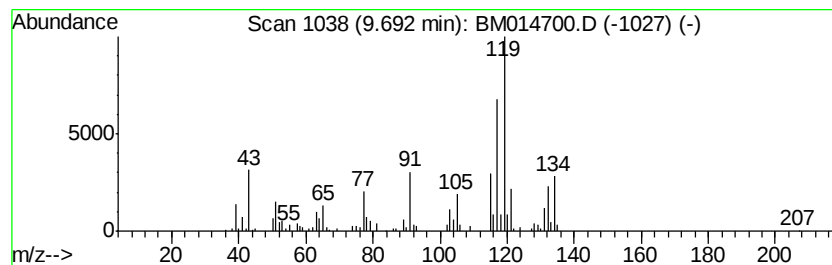
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, (1-methyl-1-propen... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	2.36 ng/ul	378220	1,4-Dichlorobenzene-d4	8.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	89
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	64
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

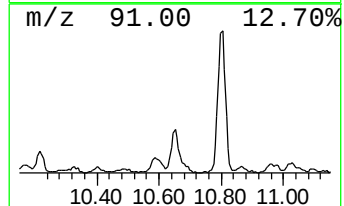
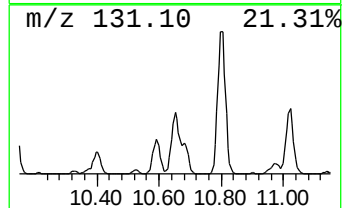
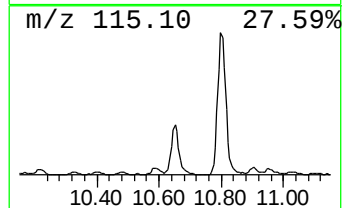
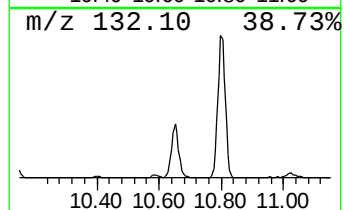
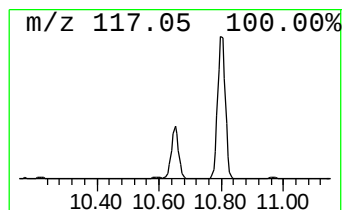
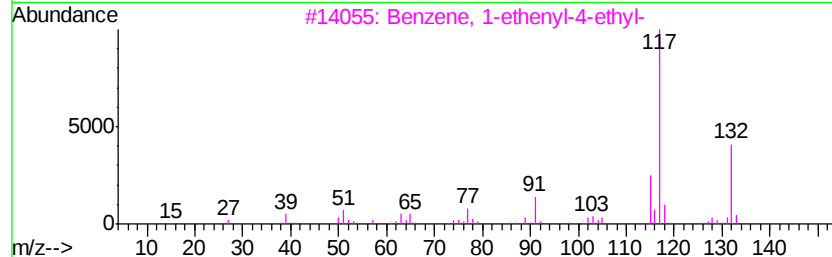
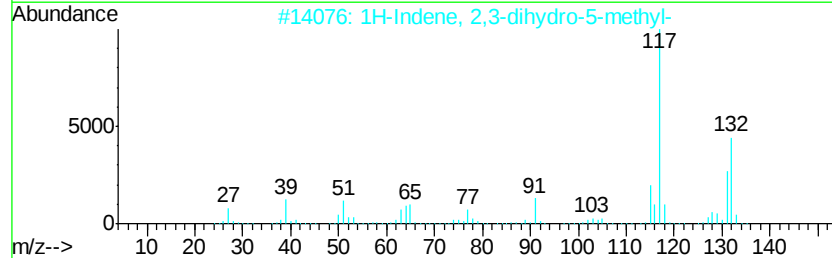
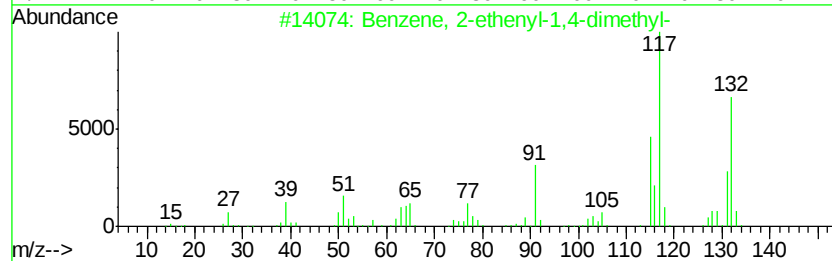
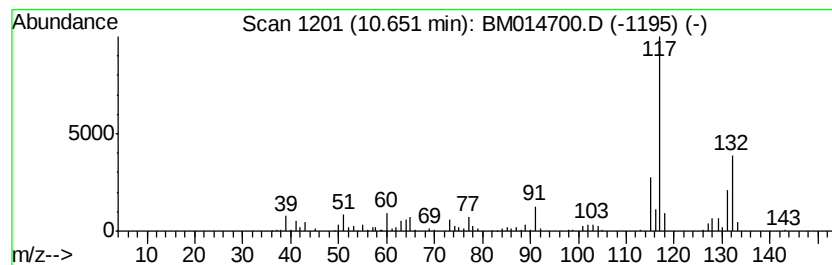
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	2.60 ng/ul	646702	Naphthalene-d8	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

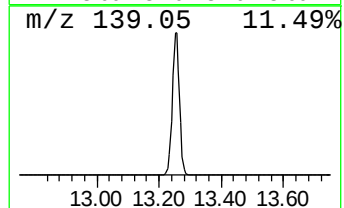
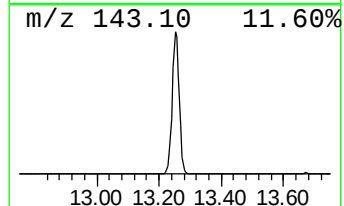
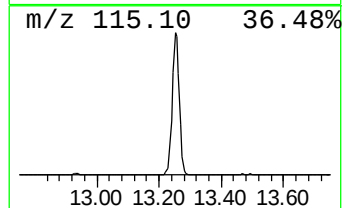
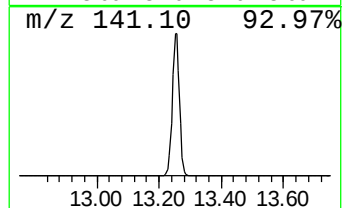
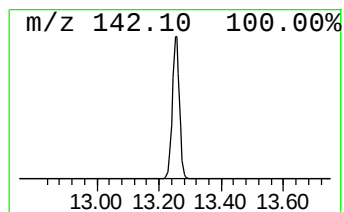
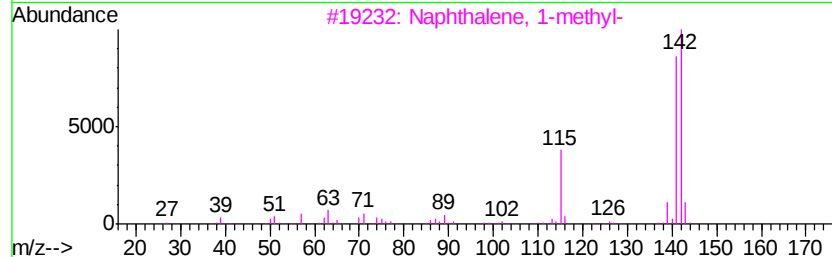
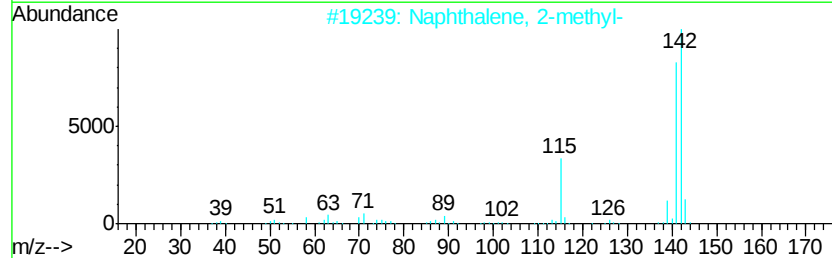
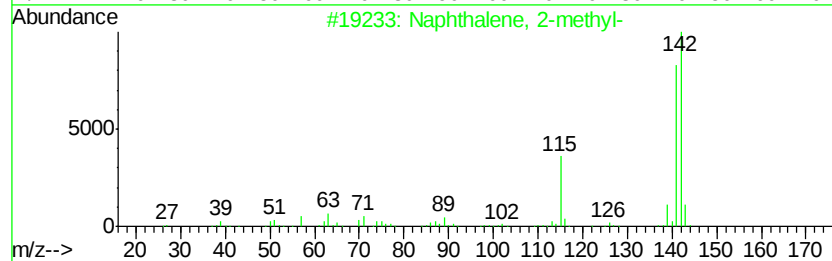
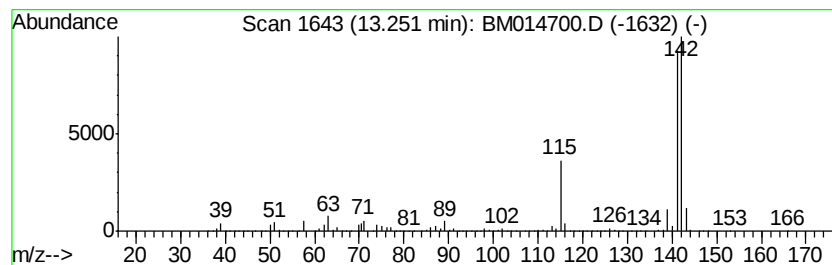
Instrument :
BNA_M
ClientSampleId :
F1A39

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.25	52.24 ng/ul	12971600	Naphthalene-d8	11.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
3		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
4		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 94
5		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

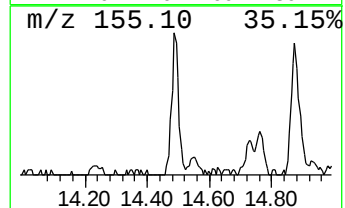
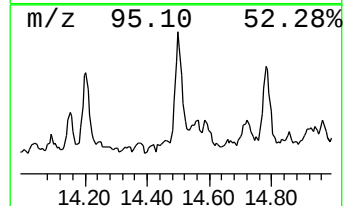
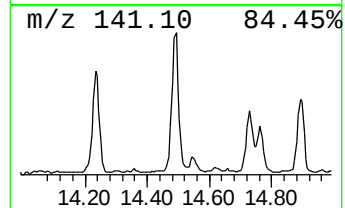
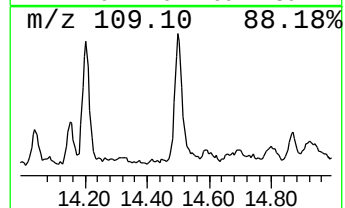
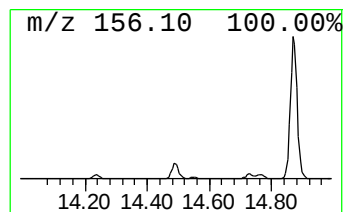
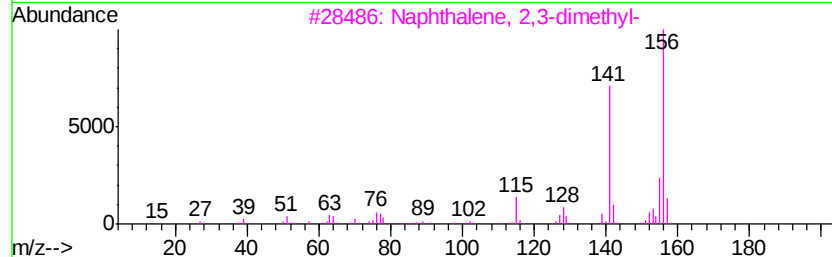
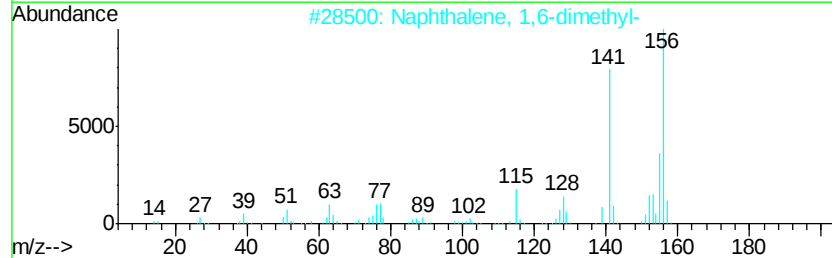
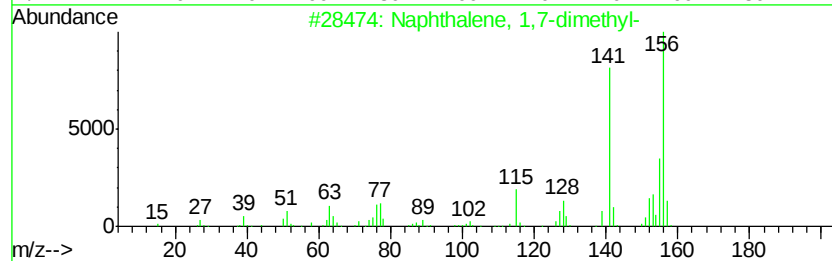
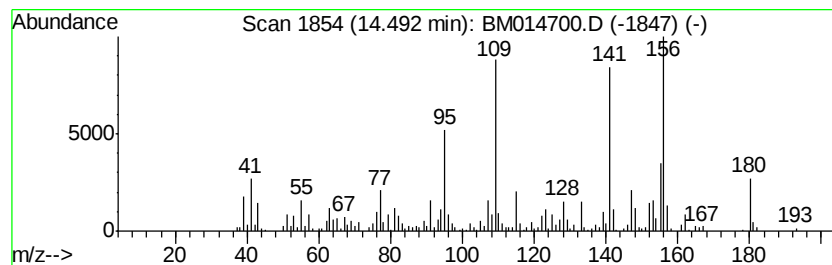
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1,7-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	2.28 ng/ul	619824	Acenaphthene-d10	15.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	91
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

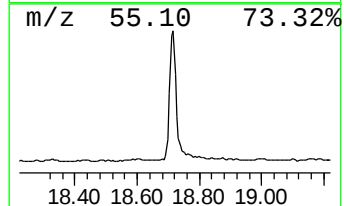
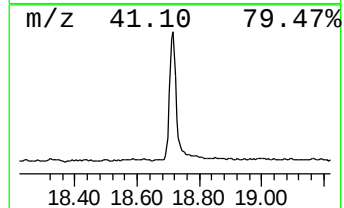
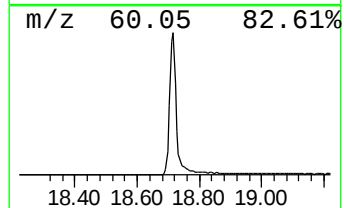
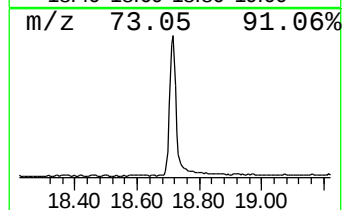
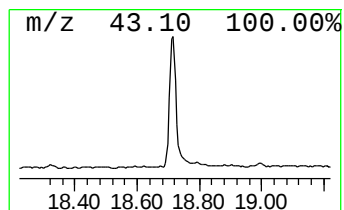
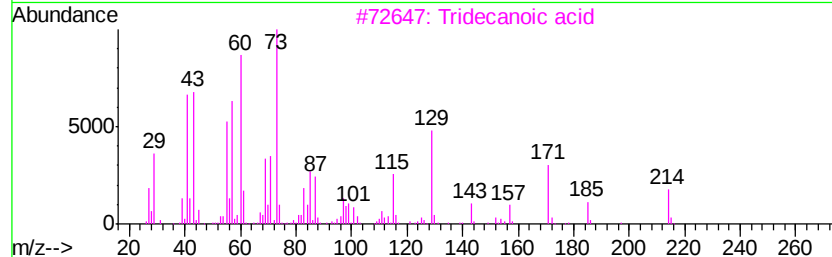
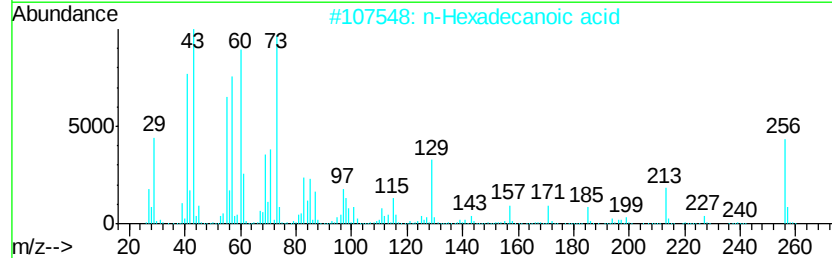
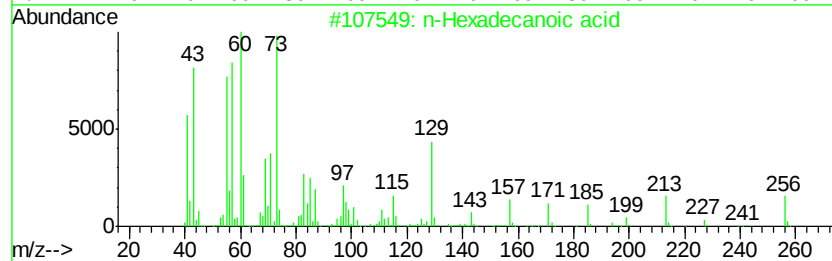
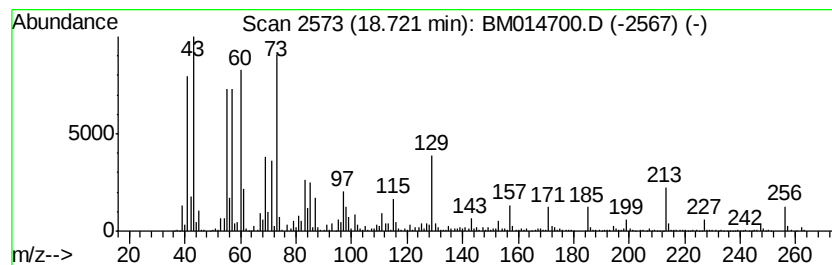
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.72	11.57 ng/ul	3246960	Phenanthrene-d10	17.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3		Tridecanoic acid	214	C13H26O2	000638-53-9	95
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	80



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

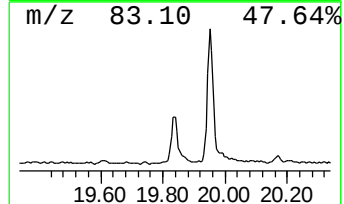
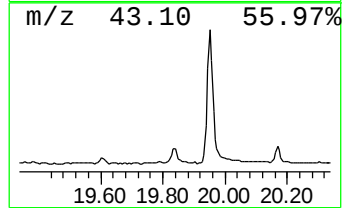
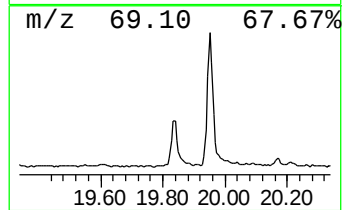
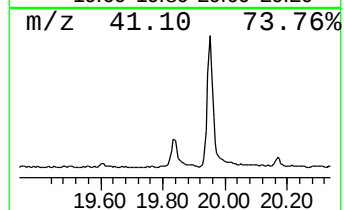
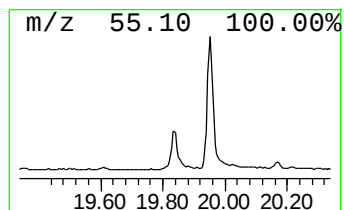
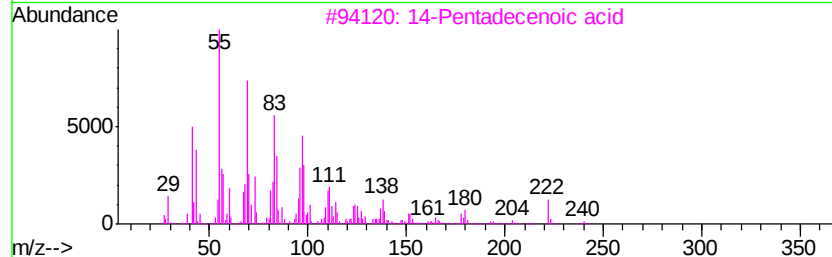
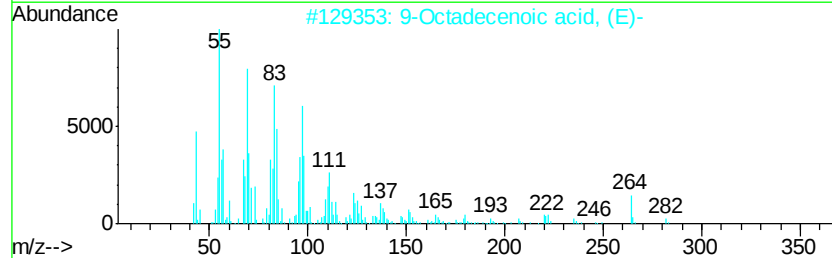
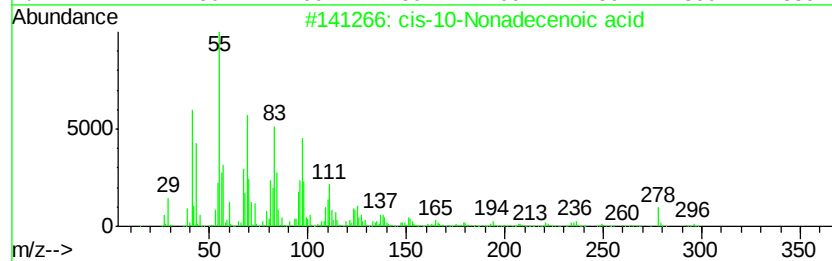
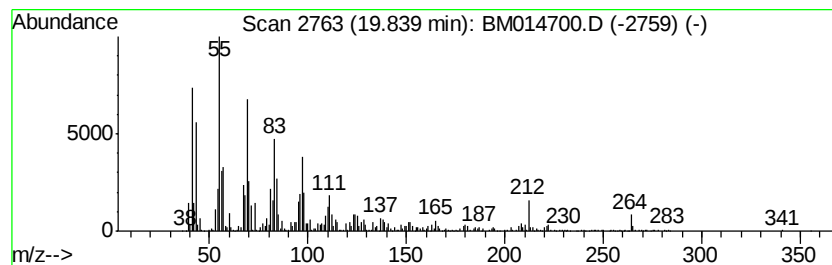
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 cis-10-Nonadecenoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.84	3.20 ng/ul	898508	Phenanthrene-d10	17.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	93
2		9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	93
3		14-Pentadecenoic acid	240	C15H28O2	017351-34-7	91
4		9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	90
5		trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

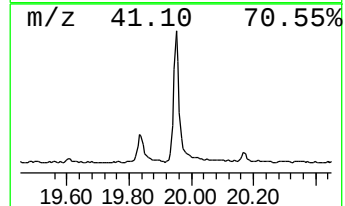
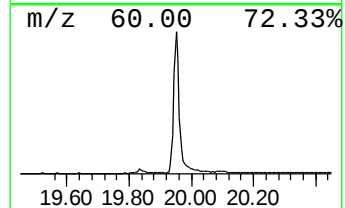
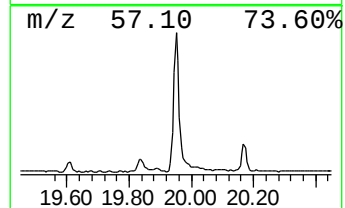
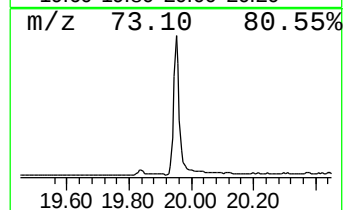
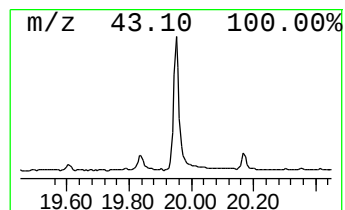
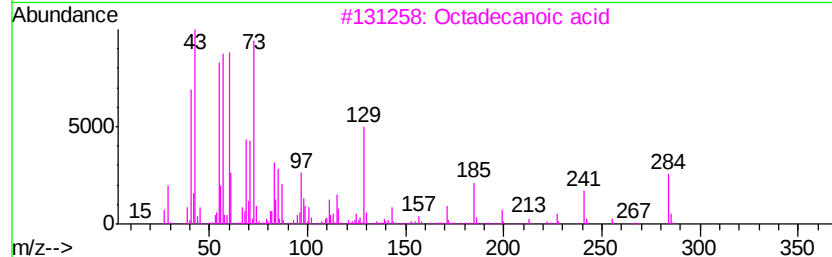
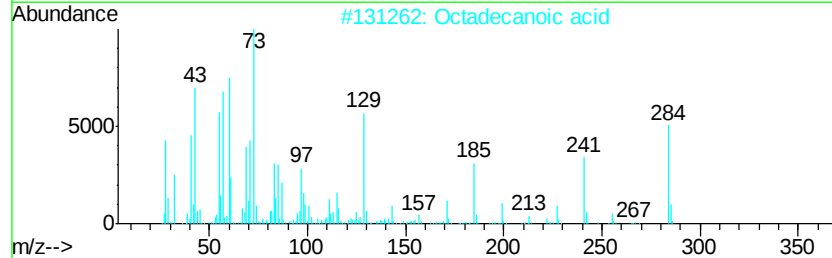
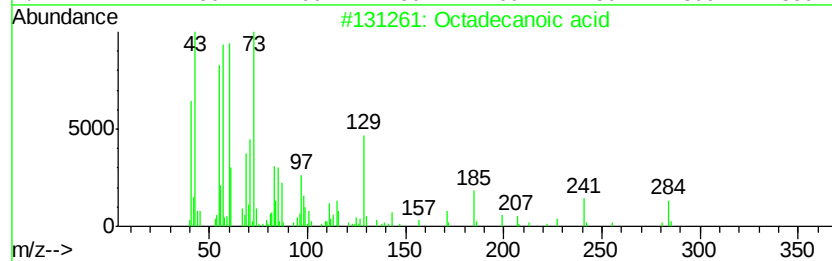
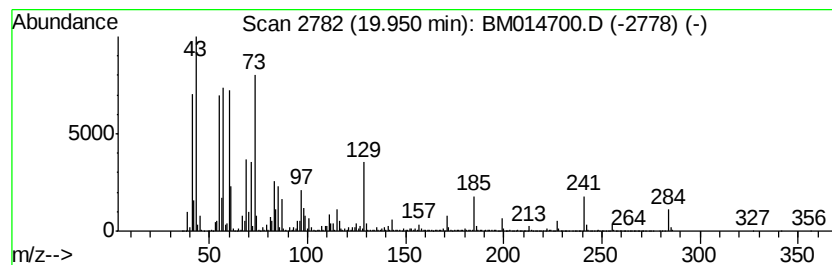
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.95	19.36 ng/ul	4373020	Chrysene-d12	21.96

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	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	98
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

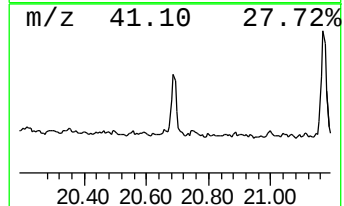
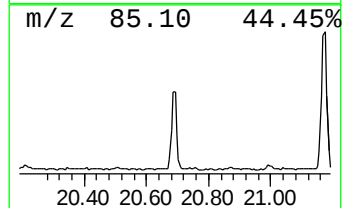
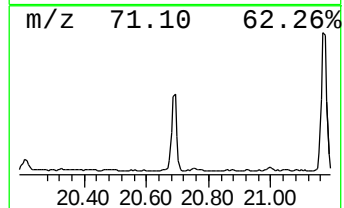
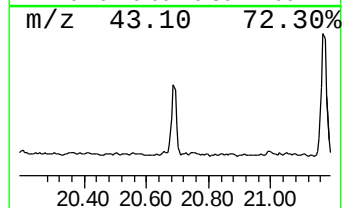
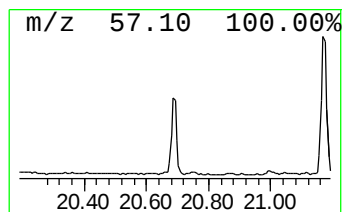
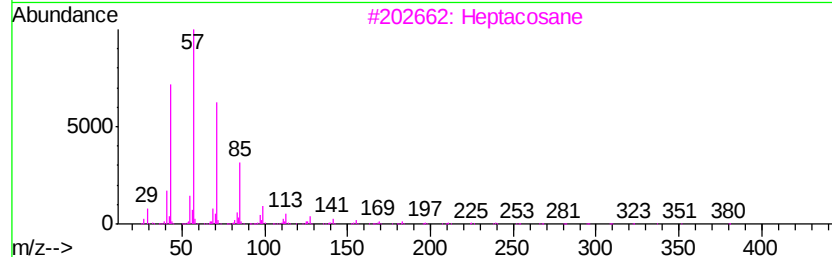
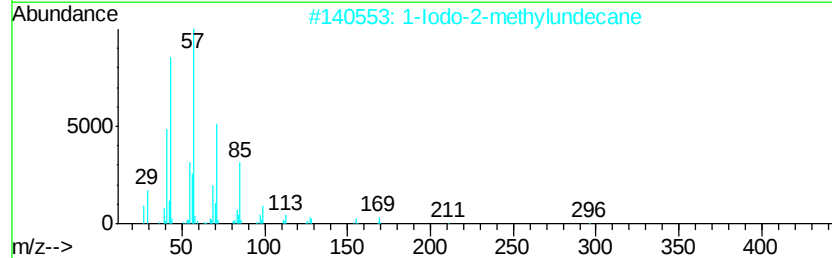
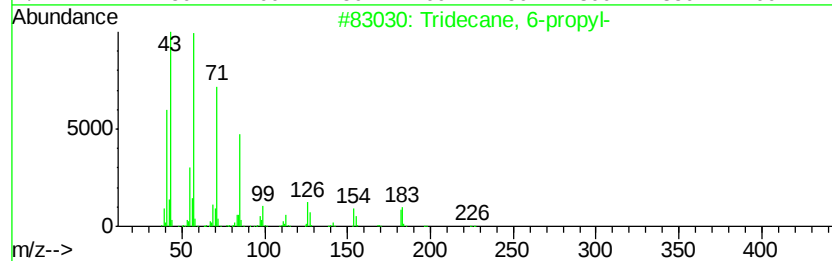
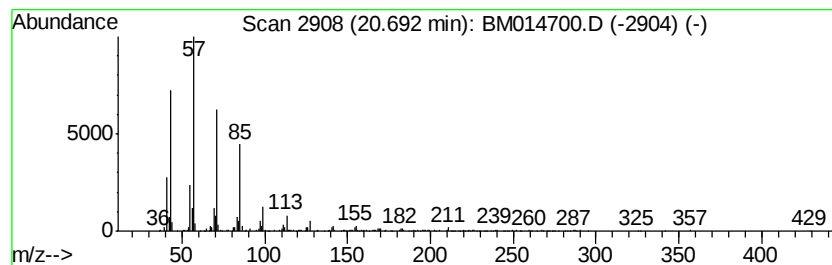
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	2.00 ng/ul	452689	Chrysene-d12	21.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 6-propyl-	226	C16H34	055045-10-8	93
2		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Eicosane	282	C20H42	000112-95-8	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

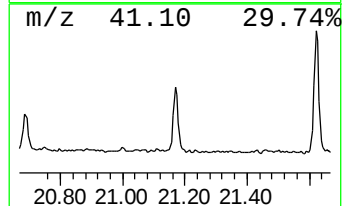
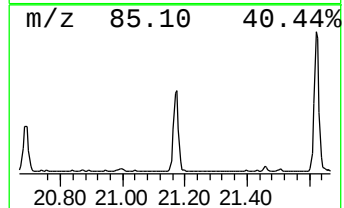
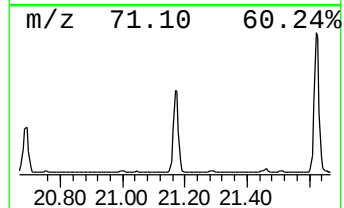
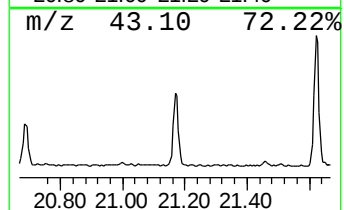
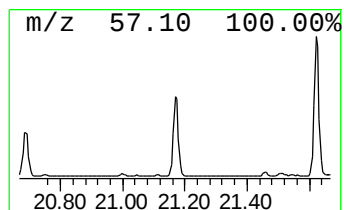
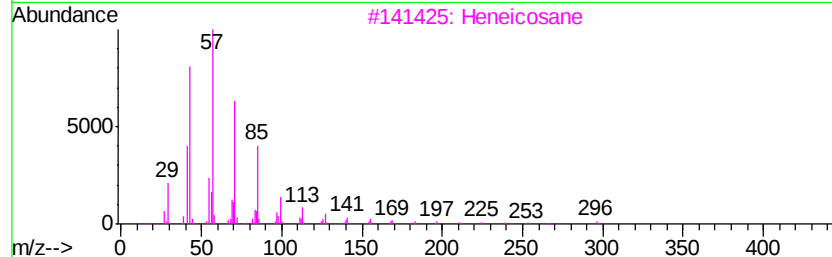
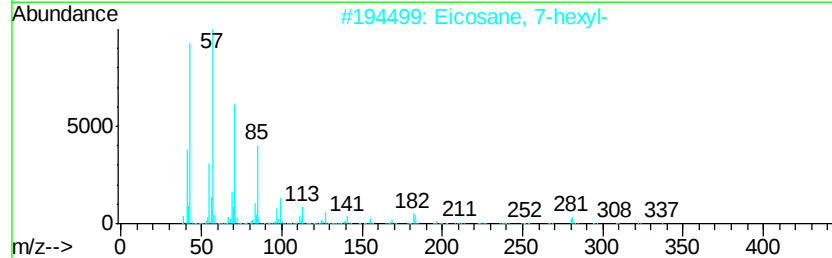
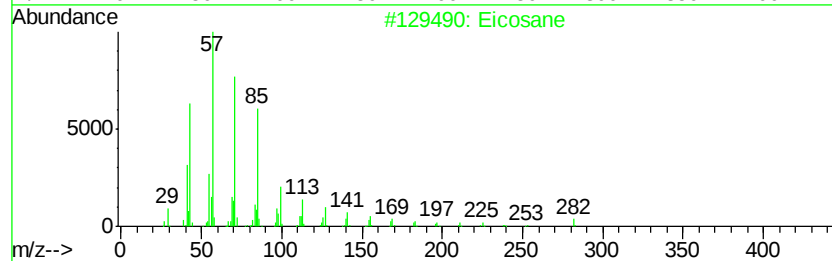
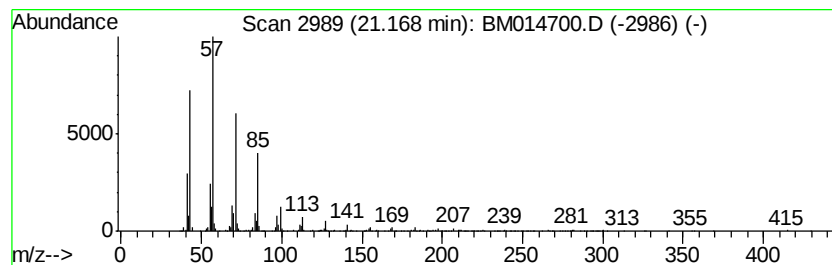
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.17	3.84 ng/ul	866389	Chrysene-d12	21.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

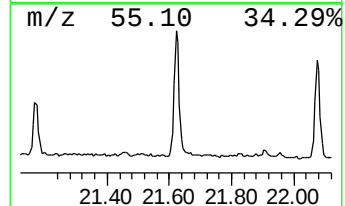
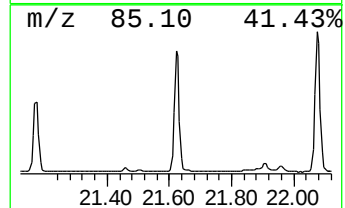
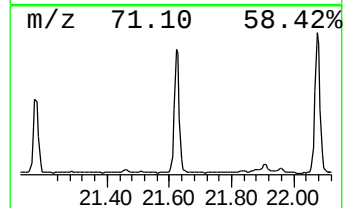
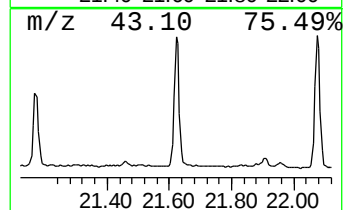
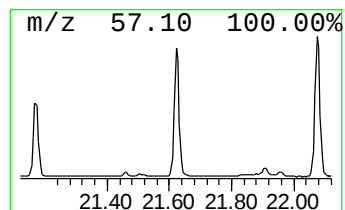
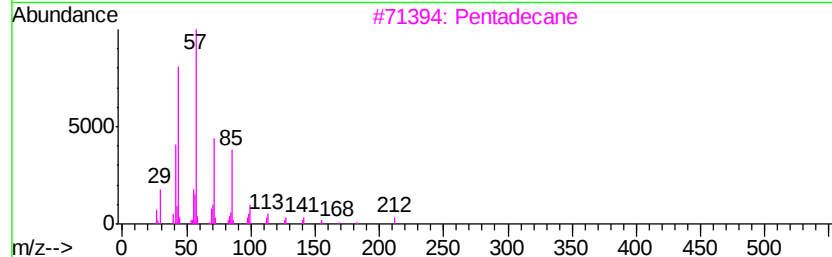
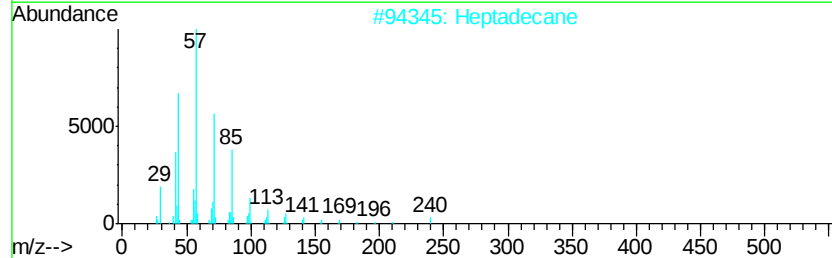
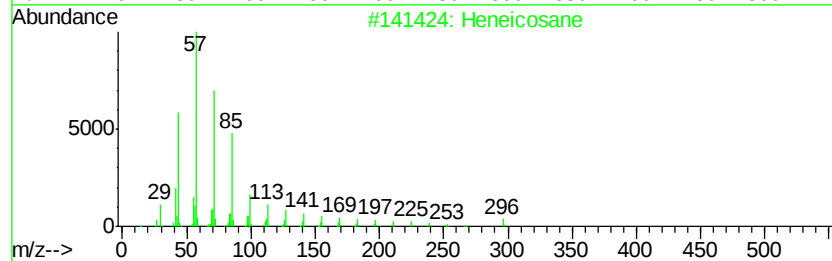
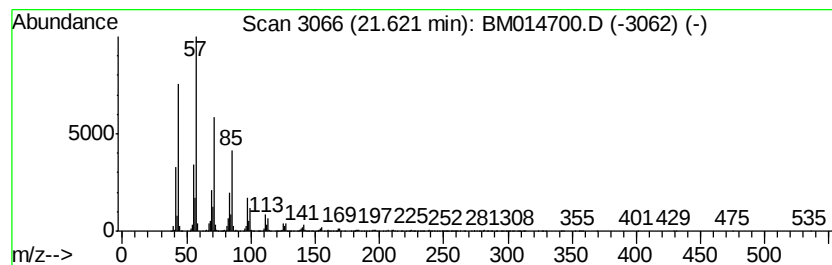
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.62	7.72 ng/ul	1744300	Chrysene-d12	21.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	94
2			Heptadecane	240	C17H36	000629-78-7	94
3			Pentadecane	212	C15H32	000629-62-9	93
4			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
5			Tritetracontane	605	C43H88	007098-21-7	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

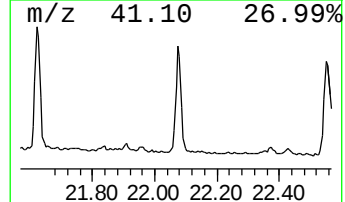
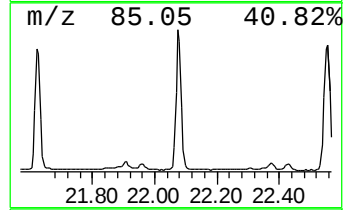
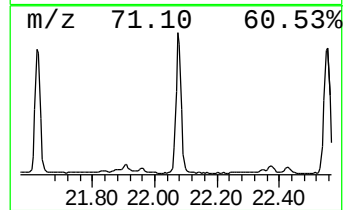
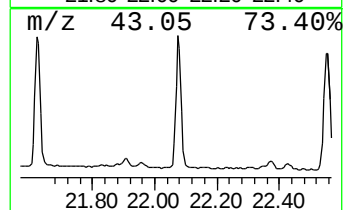
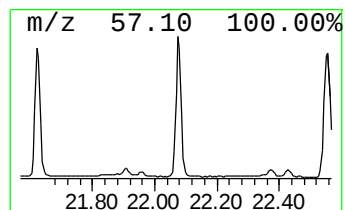
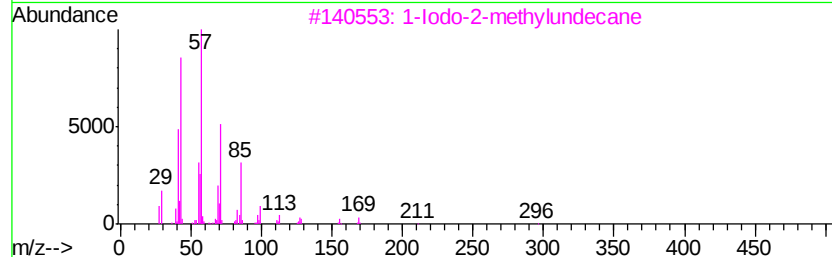
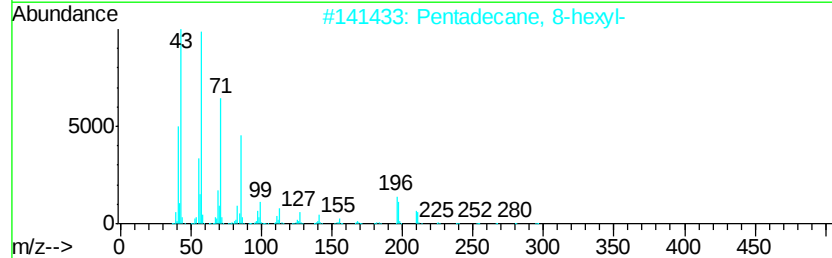
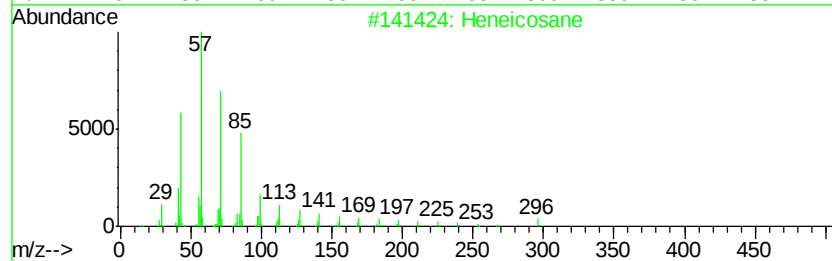
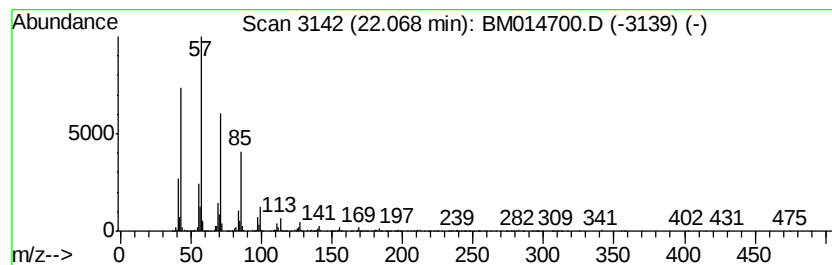
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.07	7.19 ng/ul	1625100	Chrysene-d12	21.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	95
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

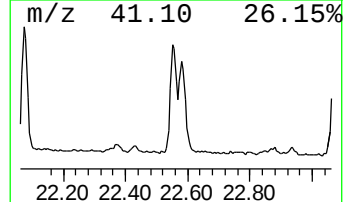
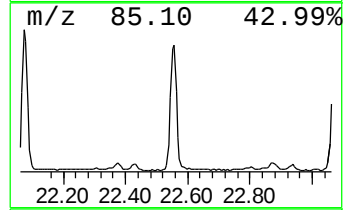
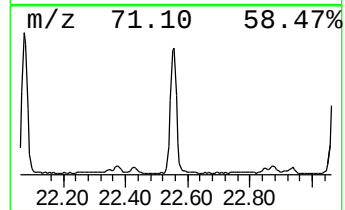
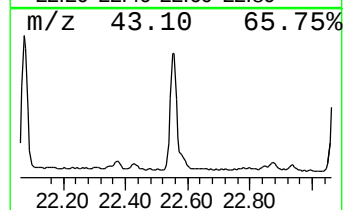
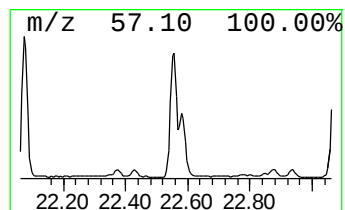
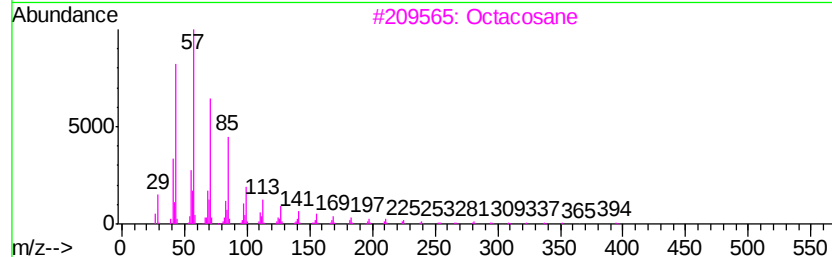
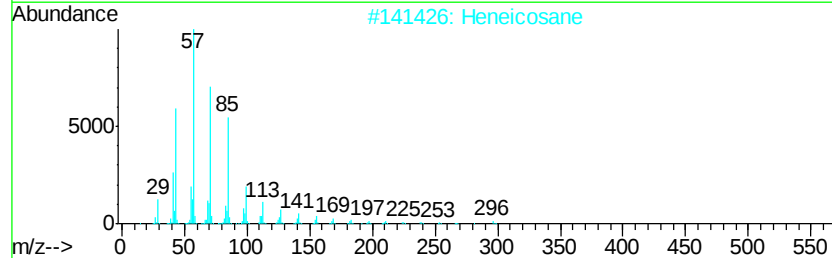
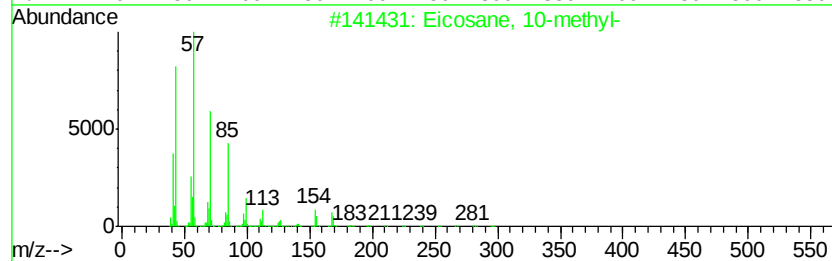
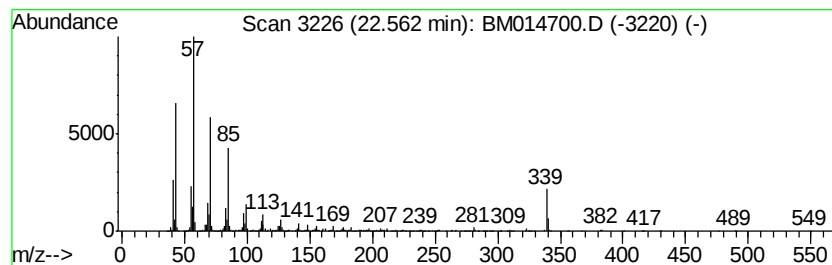
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.56	8.38 ng/ul	1893290	Chrysene-d12	21.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

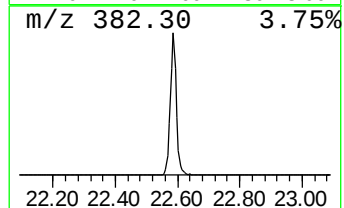
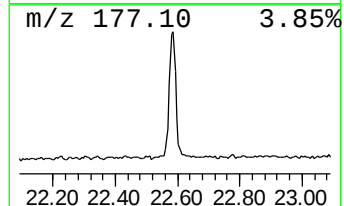
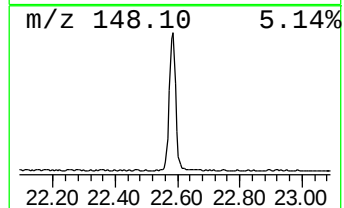
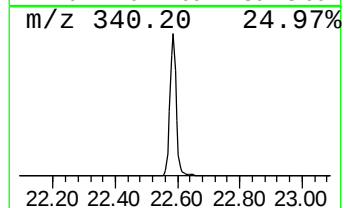
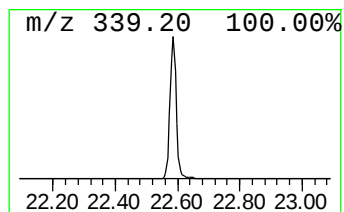
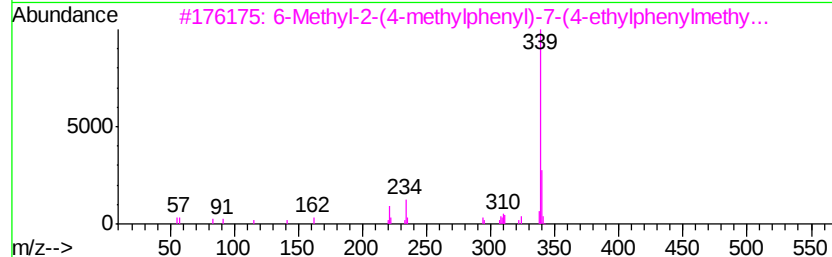
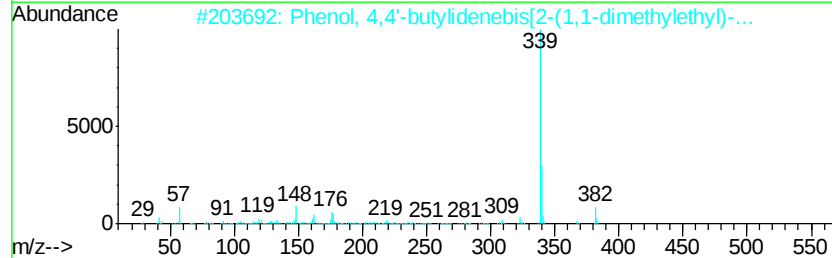
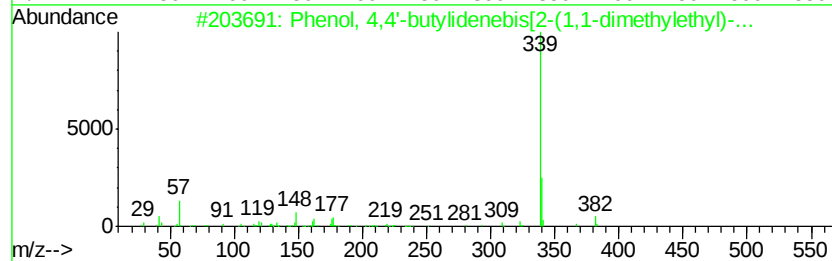
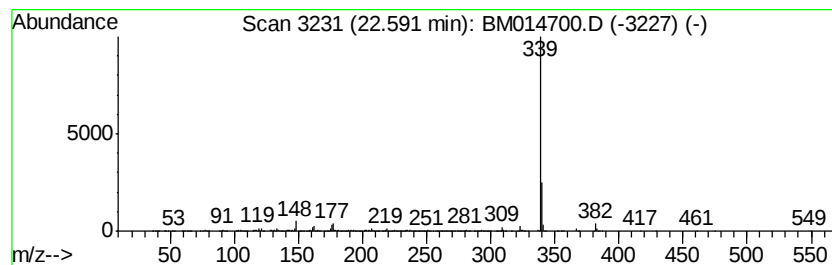
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 Phenol, 4,4'-butylidenebis[... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.59	17.12 ng/ul	3867430	Chrysene-d12	21.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-butylidenebis[2-(1,...	382	C26H38O2	000085-60-9	98
2		Phenol, 4,4'-butylidenebis[2-(1,...	382	C26H38O2	000085-60-9	90
3		6-Methyl-2-(4-methylphenyl)-7-(4...)	339	C25H25N	080193-45-9	64
4		2-(Pyrrolidin-4,6-bis(4-aminophenyl)-...	339	C21H17N5	1000078-62-7	64
5		Phenol, 4,4'-butylidenebis[2-(1,...	382	C26H38O2	000085-60-9	55



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

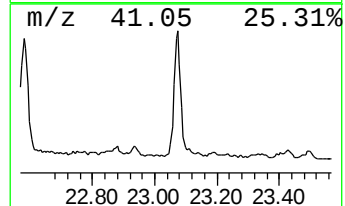
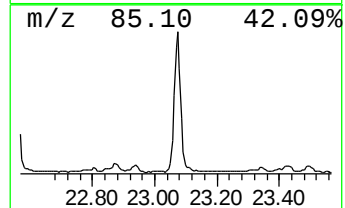
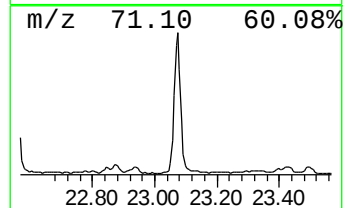
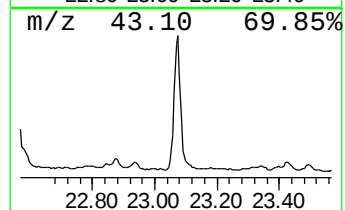
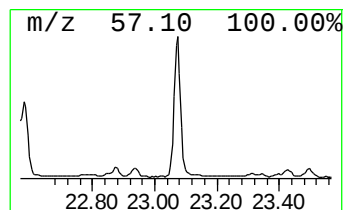
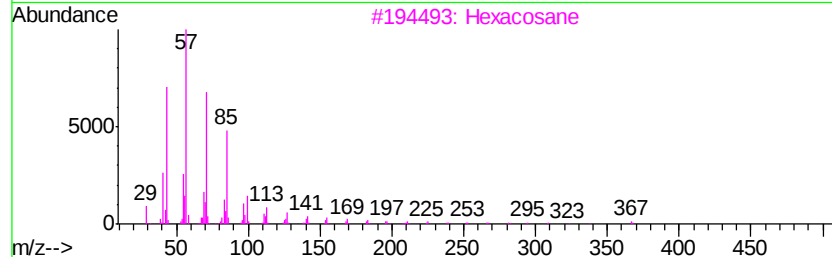
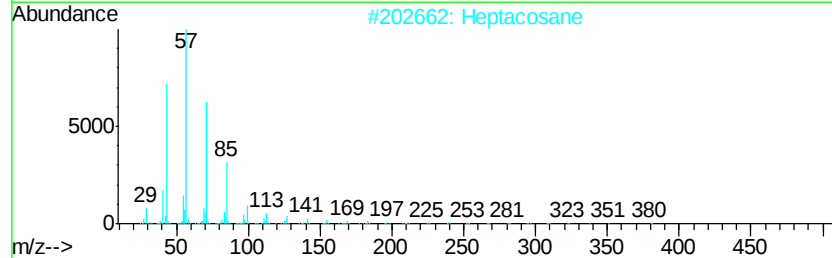
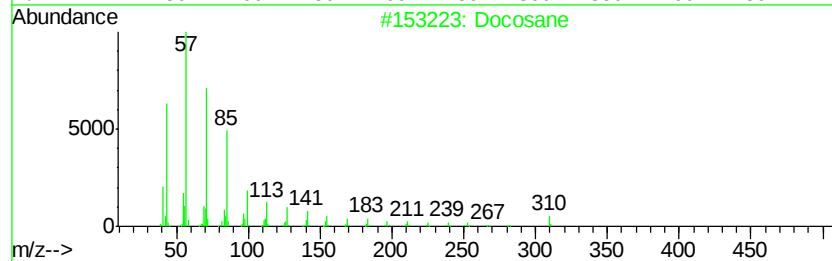
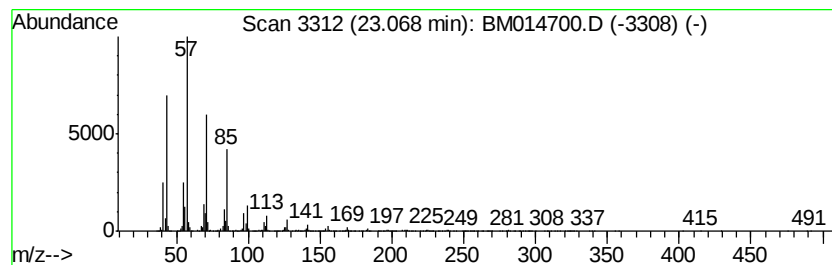
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.07	8.76 ng/ul	1978560	Chrysene-d12	21.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	97
2		Heptacosane	380	C27H56	000593-49-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

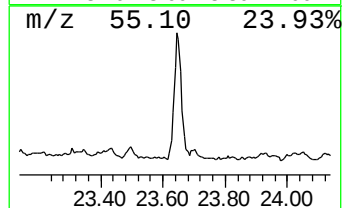
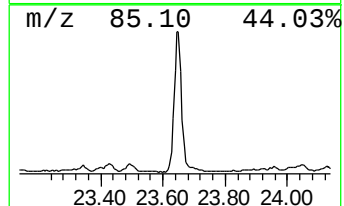
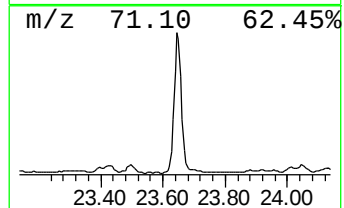
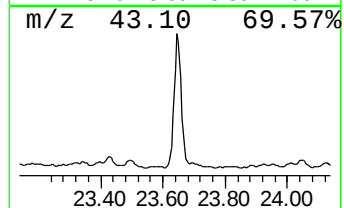
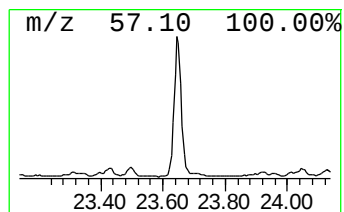
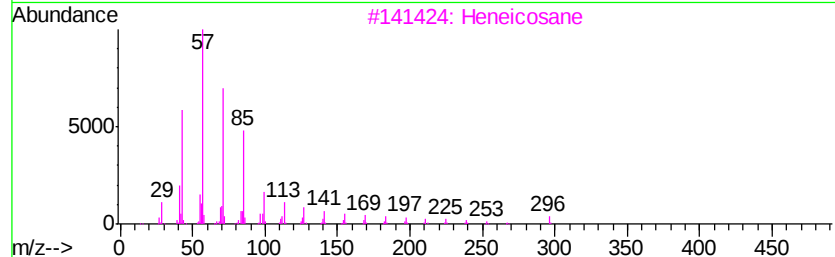
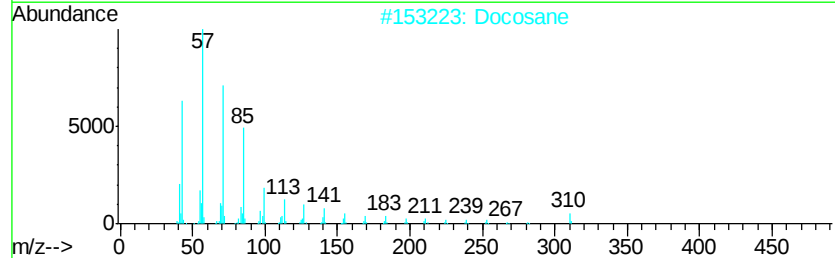
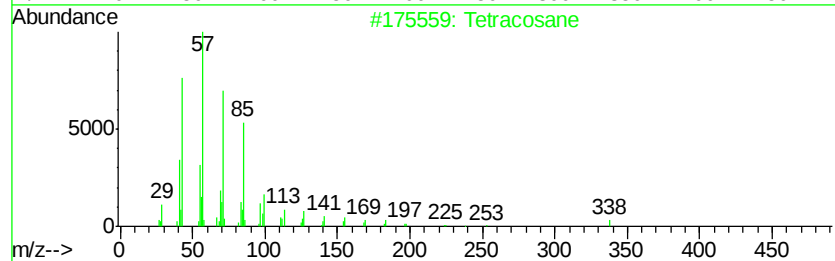
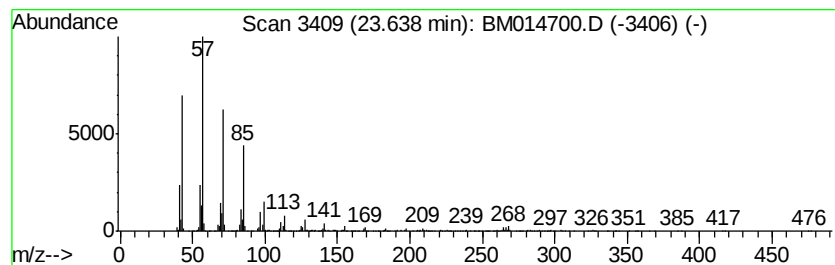
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.64	7.11 ng/ul	1532640	Perylene-d12	24.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	98
2			Docosane	310	C22H46	000629-97-0	97
3			Heneicosane	296	C21H44	000629-94-7	96
4			Nonadecane	268	C19H40	000629-92-5	94
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM041318\
Data File : BM014700.D
Acq On : 13 Apr 2018 13:43
Operator : SJ/JU
Sample : J2313-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F1A39

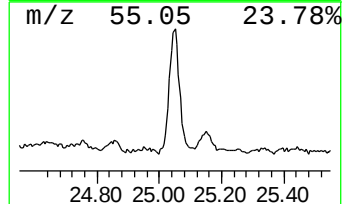
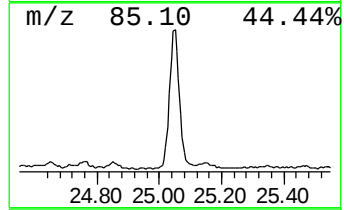
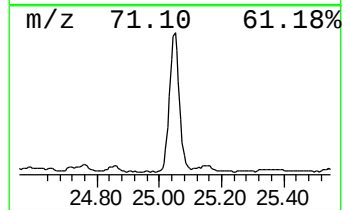
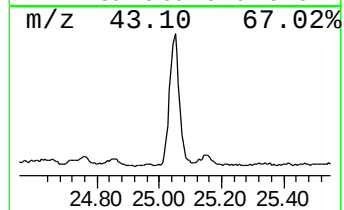
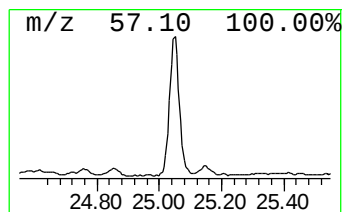
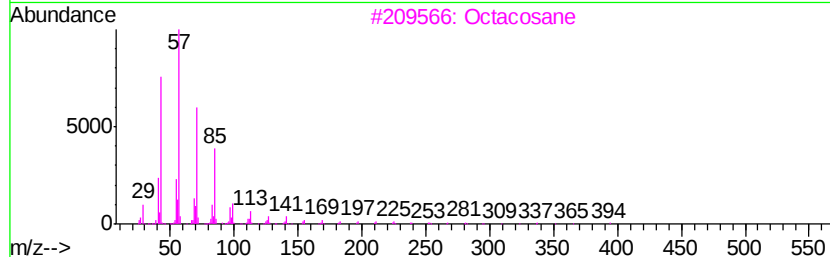
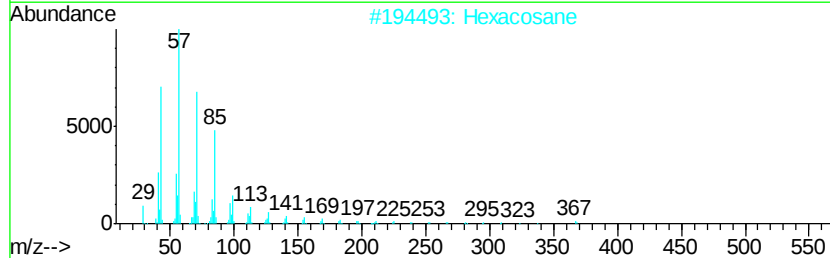
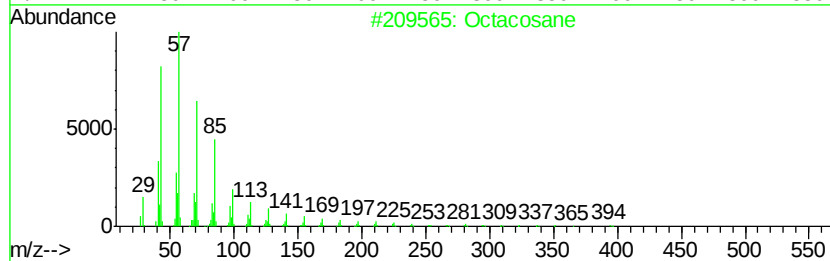
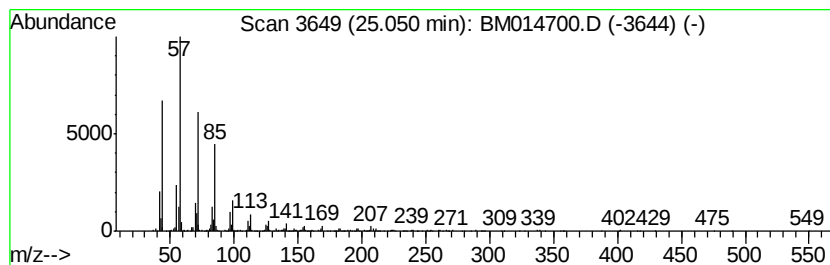
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032818MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.05	5.18 ng/ul	1116070	Perylene-d12	24.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			Octacosane	394	C28H58	000630-02-4	90
4			Tetratetracontane	619	C44H90	007098-22-8	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041318\
 Data File : BM014700.D
 Acq On : 13 Apr 2018 13:43
 Operator : SJ/JU
 Sample : J2313-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1A39

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032818MA.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
Benzene, (1-methy...	9.69	2.4	ng/ul	378220	1	8.57	3198810	20.0
Benzene, 2-etheny...	10.65	2.6	ng/ul	646702	2	11.42	4966390	20.0
Naphthalene, 1-me...	13.25	52.2	ng/ul	12971600	2	11.42	4966390	20.0
Naphthalene, 1,7-...	14.49	2.3	ng/ul	619824	3	15.17	5429150	20.0
n-Hexadecanoic acid	18.72	11.6	ng/ul	3246960	4	17.88	5611000	20.0
cis-10-Nonadeceno...	19.84	3.2	ng/ul	898508	4	17.88	5611000	20.0
Octadecanoic acid	19.95	19.4	ng/ul	4373020	5	21.96	4517310	20.0
(DEL) Alkane: Str...	20.69	2.0	ng/ul	452689	5	21.96	4517310	20.0
(DEL) Alkane: Str...	21.17	3.8	ng/ul	866389	5	21.96	4517310	20.0
(DEL) Alkane: Str...	21.62	7.7	ng/ul	1744300	5	21.96	4517310	20.0
(DEL) Alkane: Str...	22.07	7.2	ng/ul	1625100	5	21.96	4517310	20.0
(DEL) Alkane: Str...	22.56	8.4	ng/ul	1893290	5	21.96	4517310	20.0
Phenol, 4,4'-buty...	22.59	17.1	ng/ul	3867430	5	21.96	4517310	20.0
(DEL) Alkane: Str...	23.07	8.8	ng/ul	1978560	5	21.96	4517310	20.0
(DEL) Alkane: Str...	23.64	7.1	ng/ul	1532640	6	24.48	4312950	20.0
(DEL) Alkane: Str...	25.05	5.2	ng/ul	1116070	6	24.48	4312950	20.0