

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
 Data File : BM021303.D
 Acq On : 01 Jul 2019 11:11
 Operator : HP/JU
 Sample : K3590-04DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5BA3DL

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-BM061419MA.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	3.010	3	4	19	rVB	387460	451109	7.28%	0.989%
2	4.399	238	240	247	rVB2	43636	56952	0.92%	0.125%
3	6.857	654	658	663	rVB	26264	36544	0.59%	0.080%
4	6.928	665	670	684	rVB	100931	192764	3.11%	0.423%
5	7.098	694	699	705	rBV	76434	117836	1.90%	0.258%
6	7.293	726	732	739	rVB	107813	184596	2.98%	0.405%
7	7.757	804	811	819	rVB	820129	1311190	21.17%	2.875%
8	7.881	829	832	837	rVB2	10567	15477	0.25%	0.034%
9	8.463	925	931	939	rBV	132016	220972	3.57%	0.485%
10	8.922	1003	1009	1019	rVB	105500	184238	2.98%	0.404%
11	9.281	1065	1070	1075	rVB2	18998	27645	0.45%	0.061%
12	9.634	1123	1130	1138	rVB	87248	146416	2.36%	0.321%
13	10.169	1215	1221	1230	rBV	136373	245881	3.97%	0.539%
14	10.539	1277	1284	1294	rBV	1255168	2115359	34.16%	4.638%
15	10.698	1304	1311	1321	rBV	99417	196748	3.18%	0.431%
16	11.339	1415	1420	1423	rVB4	4816	6745	0.11%	0.015%
17	11.769	1488	1493	1499	rVB3	10764	17305	0.28%	0.038%
18	12.380	1594	1597	1601	rBV3	4972	7266	0.12%	0.016%
19	12.710	1647	1653	1659	rBV	98192	167234	2.70%	0.367%
20	12.775	1659	1664	1683	rVV	267690	541203	8.74%	1.187%
21	12.910	1683	1687	1694	rVV5	20891	37826	0.61%	0.083%
22	12.980	1698	1699	1707	rVB6	4834	8187	0.13%	0.018%
23	13.151	1723	1728	1729	rBV4	5875	6299	0.10%	0.014%
24	13.216	1731	1739	1749	rVV2	59844	177577	2.87%	0.389%
25	13.322	1752	1757	1762	rVV	98876	175746	2.84%	0.385%
26	13.398	1765	1770	1775	rVV	165002	296880	4.79%	0.651%
27	13.457	1775	1780	1789	rVV	161567	318861	5.15%	0.699%
28	13.551	1793	1796	1800	rVV4	7525	8420	0.14%	0.018%
29	13.592	1800	1803	1808	rVV7	7540	13242	0.21%	0.029%
30	13.710	1818	1823	1824	rBV4	12850	19182	0.31%	0.042%
31	13.810	1831	1840	1844	rBV	337087	533628	8.62%	1.170%
32	13.863	1844	1849	1853	rVV2	187323	285539	4.61%	0.626%
33	13.904	1853	1856	1863	rVB7	35810	77769	1.26%	0.171%
34	13.980	1865	1869	1873	rBV4	23818	30655	0.50%	0.067%

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35	14.092	1882	1888	1893	rBV	368387	576679	9.31%	1.264%
36	14.216	1903	1909	1914	rVB3	49513	106312	1.72%	0.233%
37	14.280	1915	1920	1925	rBV2	174057	256126	4.14%	0.562%
38	14.351	1927	1932	1934	rVV4	59805	92217	1.49%	0.202%
39	14.398	1934	1940	1949	rVB2	2104897	3150818	50.88%	6.909%
40	14.627	1974	1979	1985	rBV3	102168	204527	3.30%	0.448%
41	14.698	1988	1991	1993	rBV3	37794	53254	0.86%	0.117%
42	14.786	2002	2006	2010	rVB2	81270	107471	1.74%	0.236%
43	14.839	2012	2015	2020	rVV6	45109	75093	1.21%	0.165%
44	14.904	2021	2026	2032	rVB4	141095	244509	3.95%	0.536%
45	14.974	2035	2038	2041	rVB2	58650	63812	1.03%	0.140%
46	15.033	2041	2048	2053	rBV7	84868	190545	3.08%	0.418%
47	15.239	2079	2083	2088	rBV2	373467	499838	8.07%	1.096%
48	15.345	2098	2101	2105	rBV4	73781	120298	1.94%	0.264%
49	15.392	2105	2109	2114	rVV	570302	754660	12.19%	1.655%
50	15.527	2129	2132	2138	rBV5	71344	117671	1.90%	0.258%
51	15.645	2145	2152	2156	rBV	479934	866975	14.00%	1.901%
52	15.739	2165	2168	2170	rBV3	110520	145759	2.35%	0.320%
53	15.798	2175	2178	2183	rVV5	80457	143441	2.32%	0.315%
54	15.857	2185	2188	2191	rVB2	127811	146444	2.36%	0.321%
55	15.892	2191	2194	2200	rBV3	128769	198282	3.20%	0.435%
56	15.998	2209	2212	2220	rVB4	183320	291801	4.71%	0.640%
57	16.121	2226	2233	2240	rBV	1478590	3059500	49.41%	6.708%
58	16.257	2253	2256	2260	rBV4	107232	134655	2.17%	0.295%
59	16.445	2285	2288	2291	rBV2	147835	203671	3.29%	0.447%
60	16.798	2343	2348	2354	rBV	4366788	6192320	100.00%	13.578%
61	16.892	2361	2364	2368	rVB	503456	585157	9.45%	1.283%
62	16.945	2369	2373	2377	rVV2	641265	844056	13.63%	1.851%
63	17.151	2402	2408	2413	rBV2	2740589	3907934	63.11%	8.569%
64	17.245	2421	2424	2431	rVB2	642170	935542	15.11%	2.051%
65	17.545	2473	2475	2480	rBV4	195536	275489	4.45%	0.604%
66	17.633	2486	2490	2495	rVB	666167	811798	13.11%	1.780%
67	18.033	2556	2558	2562	rVB3	155301	156382	2.53%	0.343%
68	18.327	2604	2608	2613	rVB	534684	678265	10.95%	1.487%
69	18.862	2696	2699	2708	rVB3	287051	458438	7.40%	1.005%
70	18.962	2713	2716	2719	rVB	288618	310676	5.02%	0.681%
71	19.092	2736	2738	2741	rBV2	129960	134874	2.18%	0.296%

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72	19.203	2755	2757	2764	rVB	229465	320802	5.18%	0.703%
73	19.545	2811	2815	2817	rBV	767740	959766	15.50%	2.104%
74	19.574	2817	2820	2826	rVV2	478771	826668	13.35%	1.813%
75	19.645	2829	2832	2838	rVB	362172	575035	9.29%	1.261%
76	20.368	2952	2955	2958	rBV2	147203	187432	3.03%	0.411%
77	20.533	2980	2983	2987	rBV3	399220	507921	8.20%	1.114%
78	21.339	3115	3120	3124	rBV	2284789	3021530	48.79%	6.625%
79	23.462	3477	3481	3487	rVB	500620	833636	13.46%	1.828%
80	23.609	3500	3506	3514	rVB	1606055	3045567	49.18%	6.678%

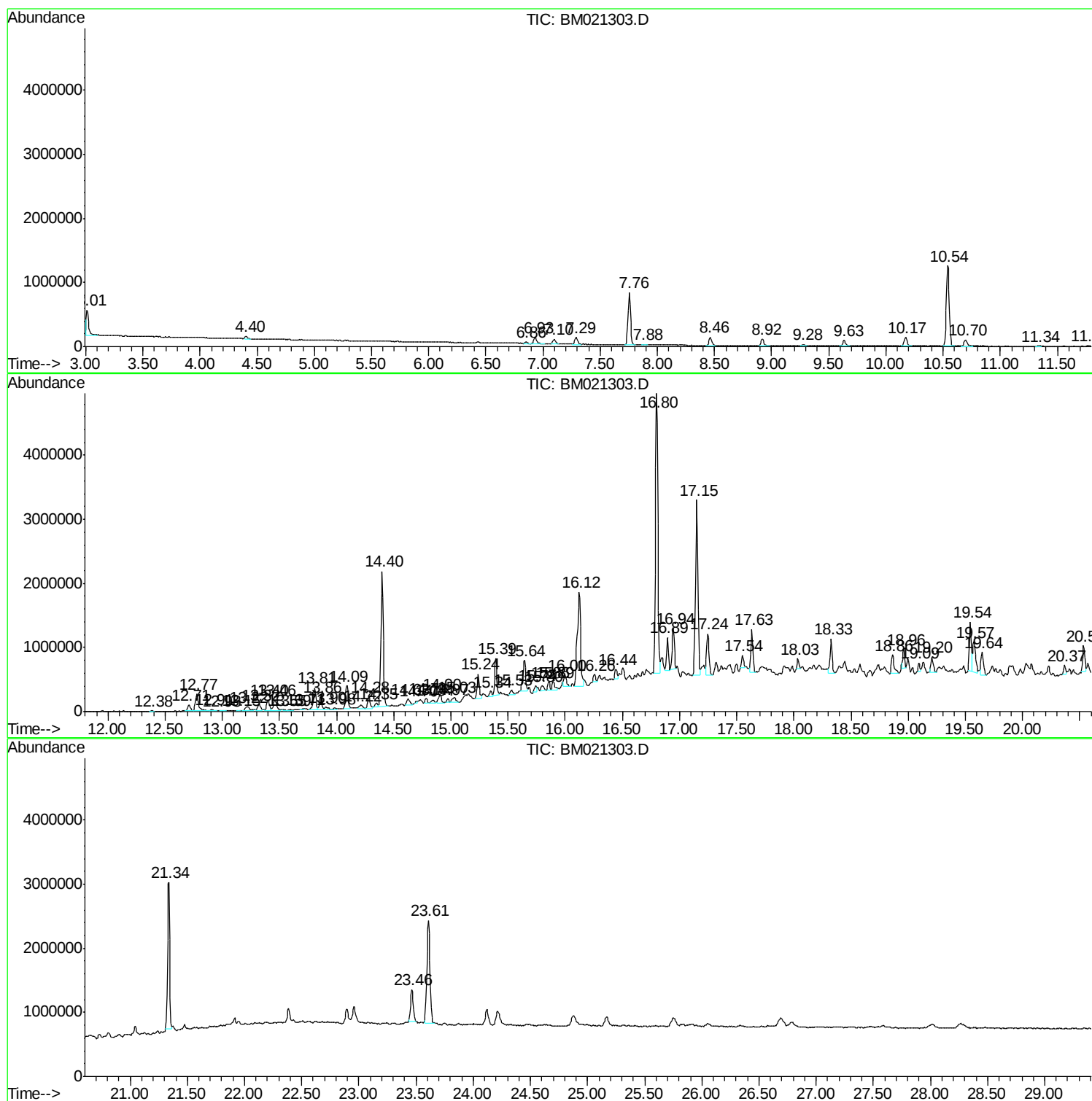
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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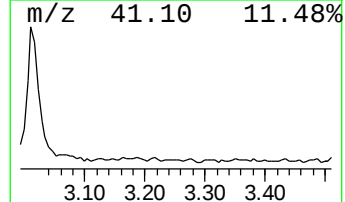
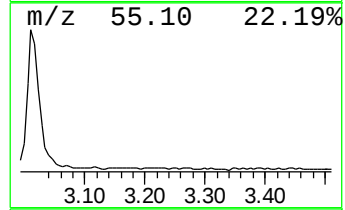
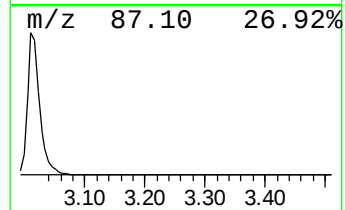
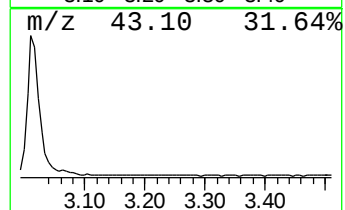
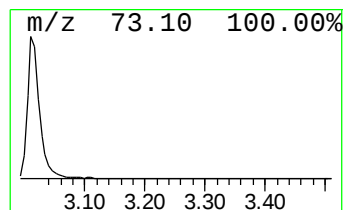
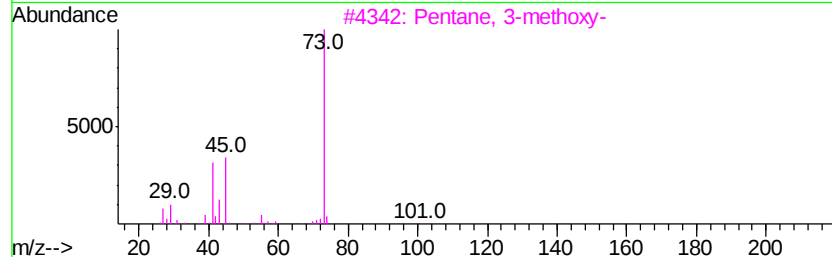
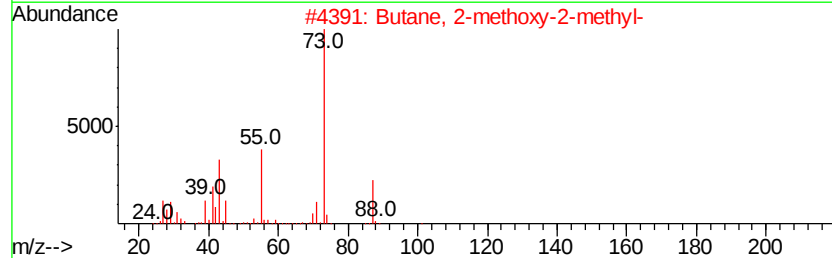
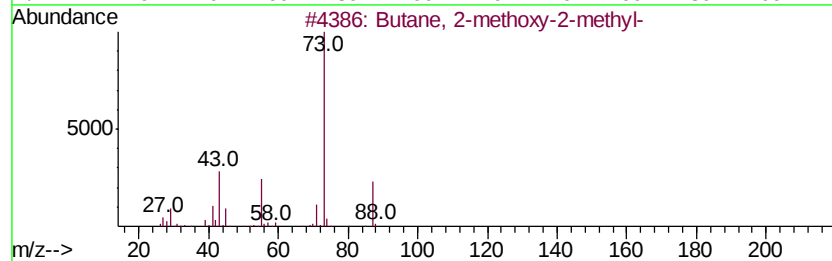
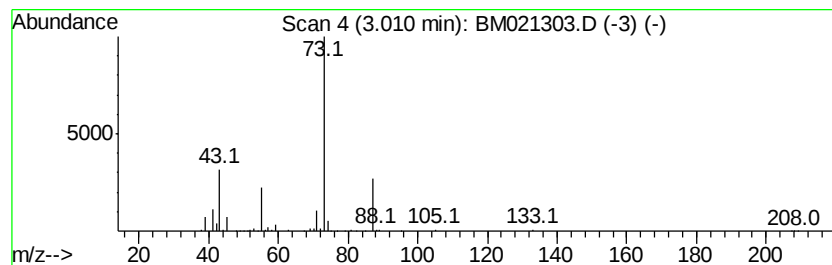
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.01	6.88 ng/ul	451109	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	45
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	38
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	16



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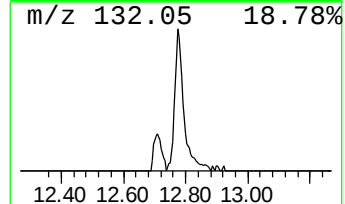
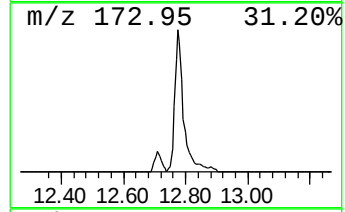
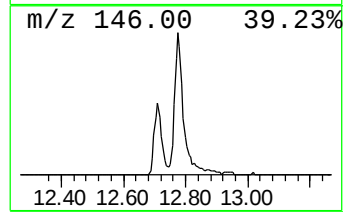
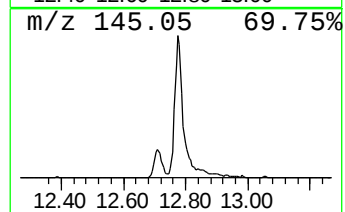
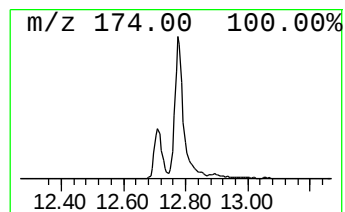
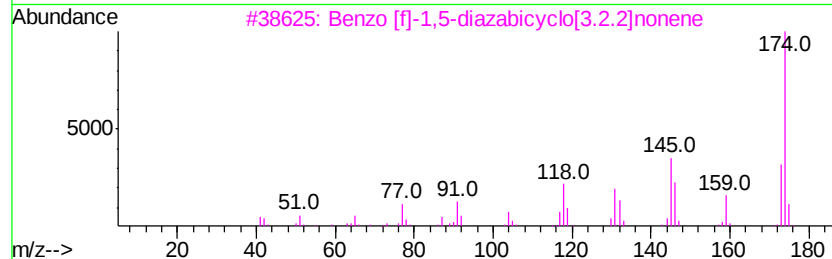
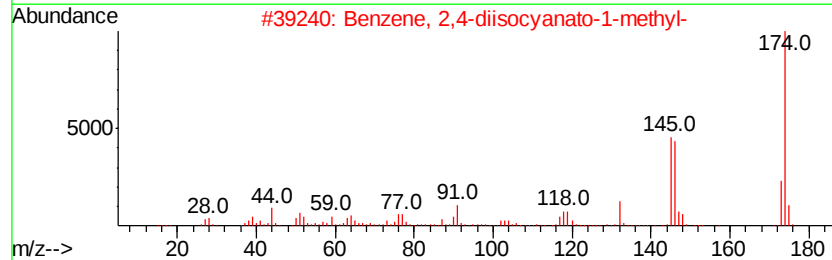
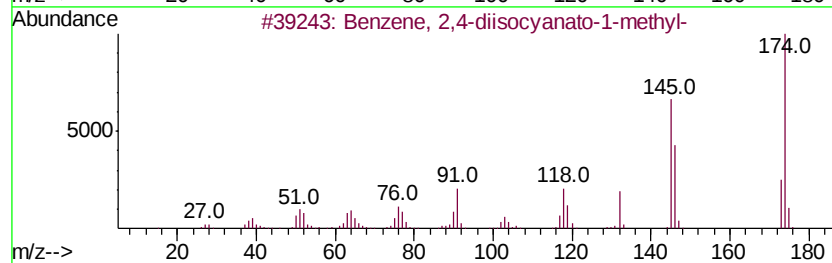
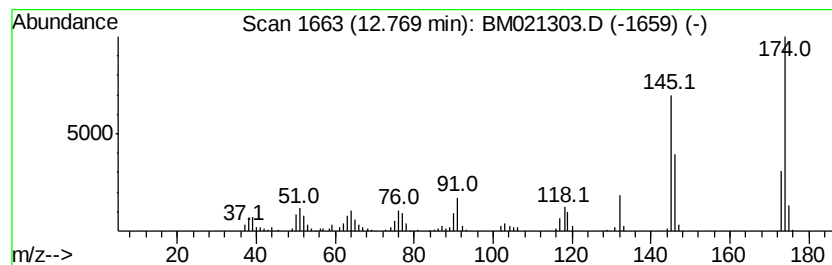
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 2,4-diisocyanato-1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	3.44 ng/ul	541203	Acenaphthene-d10	14.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2,4-diisocyanato-1-methyl-	174	C9H6N2O2	000584-84-9	93
2		Benzene, 2,4-diisocyanato-1-methyl-	174	C9H6N2O2	000584-84-9	93
3		Benzo [f]-1,5-diazabicyclo[3.2.2]nonene	174	C11H14N2	007092-76-4	72
4		Benzene, 1,3-diisocyanato-2-methyl-	174	C9H6N2O2	000091-08-7	58
5		Benzene, 2,4-diisocyanato-1-methyl-	174	C9H6N2O2	000584-84-9	52



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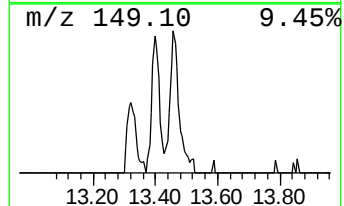
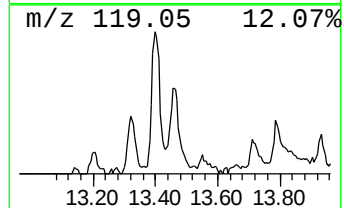
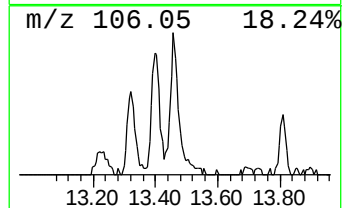
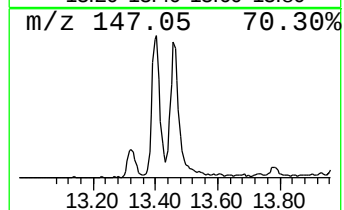
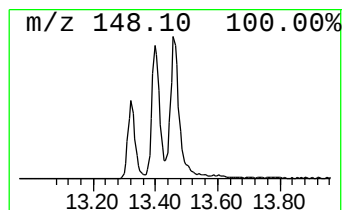
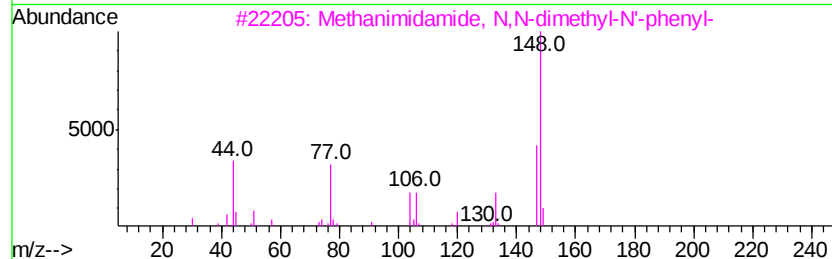
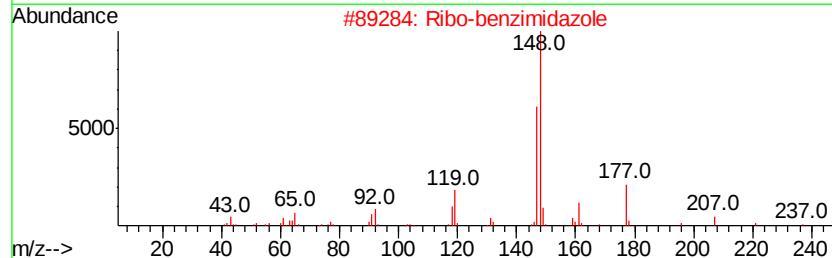
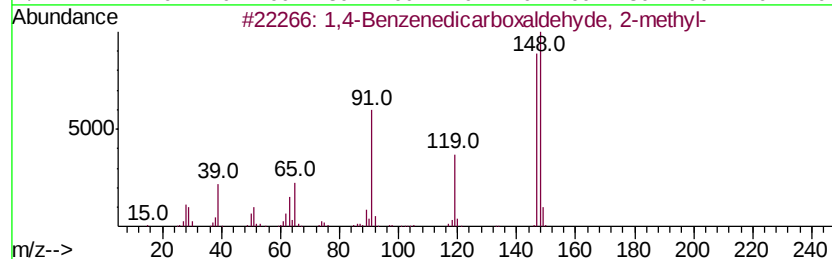
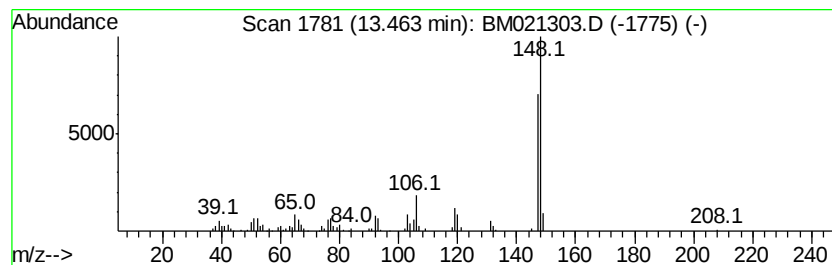
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,4-Benzenedicarboxaldehyde... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	2.02 ng/ul	318861	Acenaphthene-d10	14.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Benzenedicarboxaldehyde, 2-m...	148	C9H8O2	027587-17-3	64
2		Ribo-benzimidazole	252	C12H16N2O4	1000128-23-5	59
3		Methanimidamide, N,N-dimethyl-N'...	148	C9H12N2	001783-25-1	59
4		3H-Indazol-3-one, 1,2-dihydro-2-...	148	C8H8N2O	001848-40-4	53
5		1H-Benzotriazol-5-amine, 1-methyl-	148	C7H8N4	027799-83-3	47



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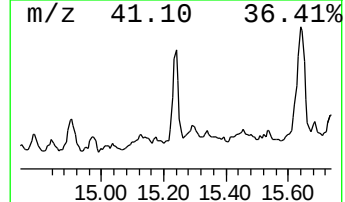
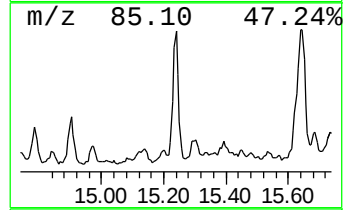
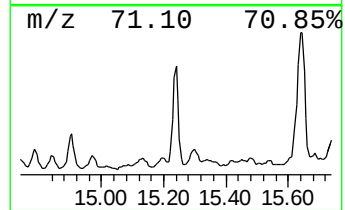
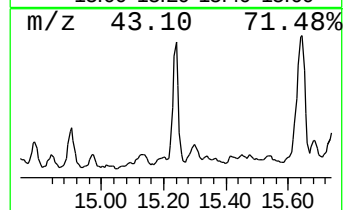
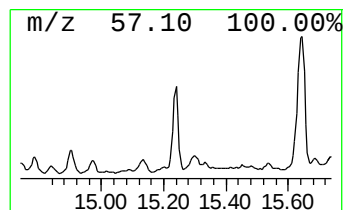
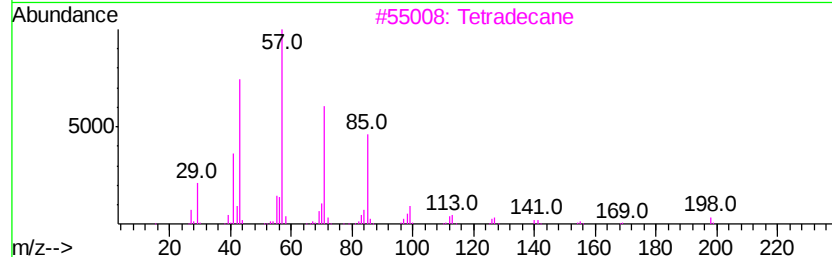
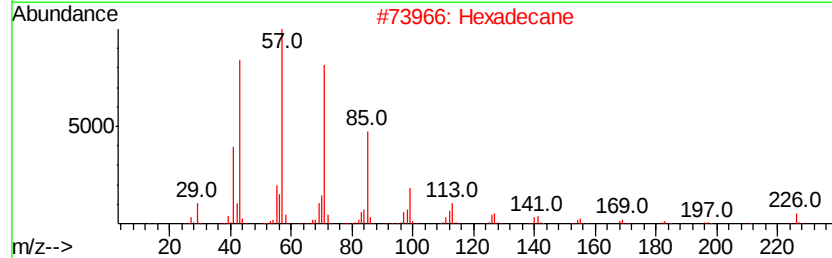
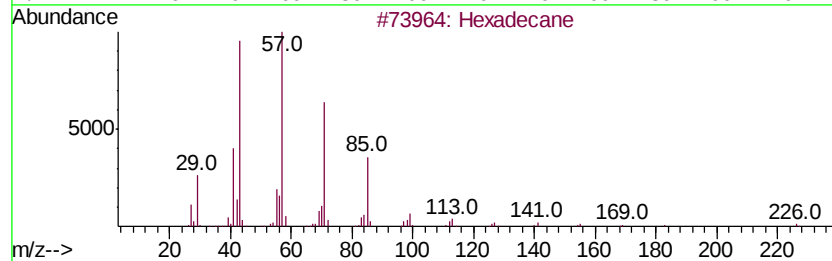
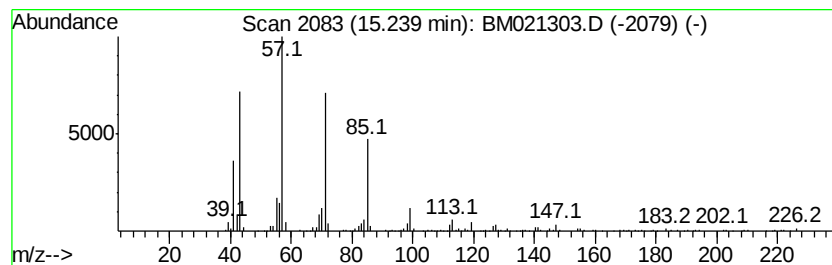
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	3.17 ng/ul	499838	Acenaphthene-d10	14.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Hexadecane	226	C16H34	000544-76-3	89
3		Tetradecane	198	C14H30	000629-59-4	87
4		Heneicosane	296	C21H44	000629-94-7	86
5		Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5BA3DL

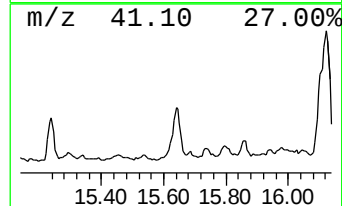
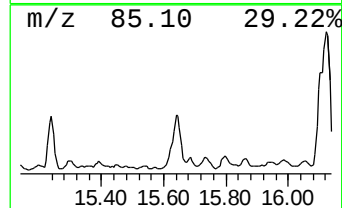
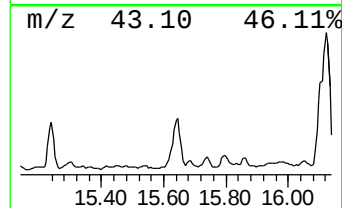
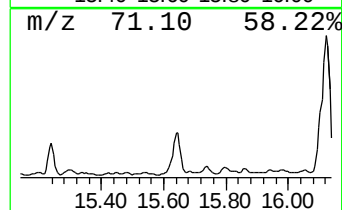
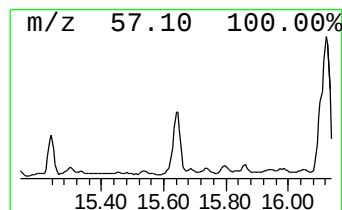
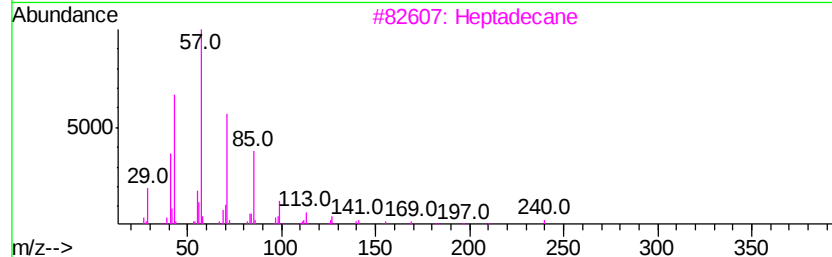
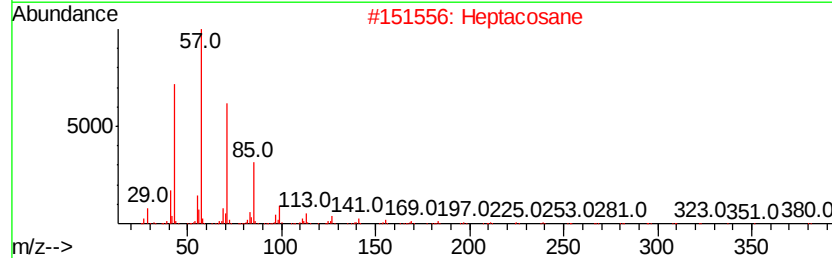
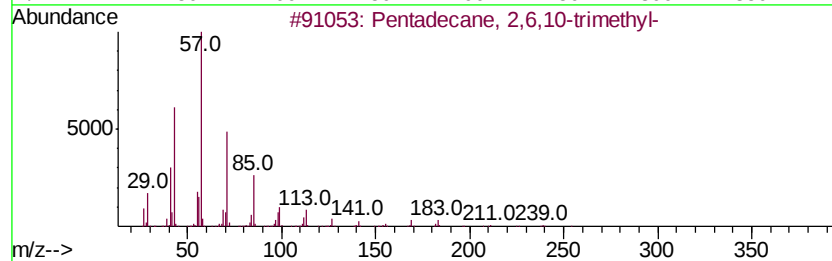
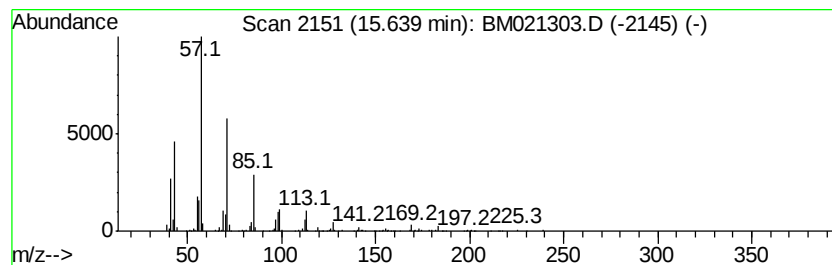
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	5.50 ng/ul	866975	Acenaphthene-d10	14.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	91
2		Heptacosane	380	C27H56	000593-49-7	80
3		Heptadecane	240	C17H36	000629-78-7	78
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
5		Pentacosane	352	C25H52	000629-99-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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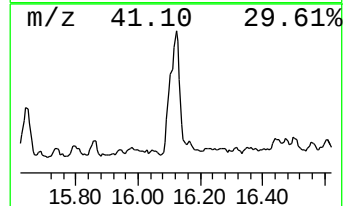
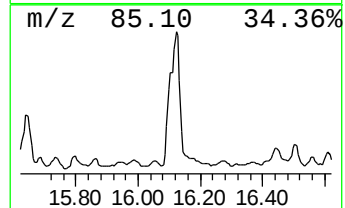
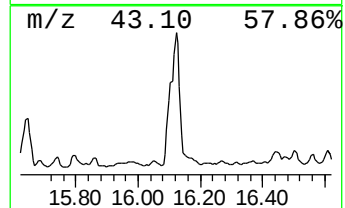
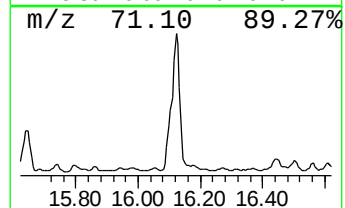
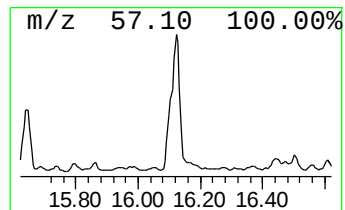
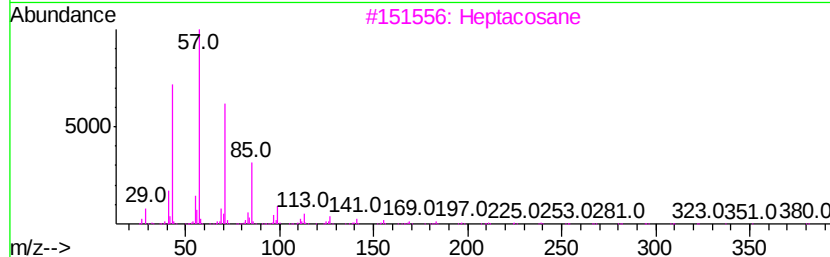
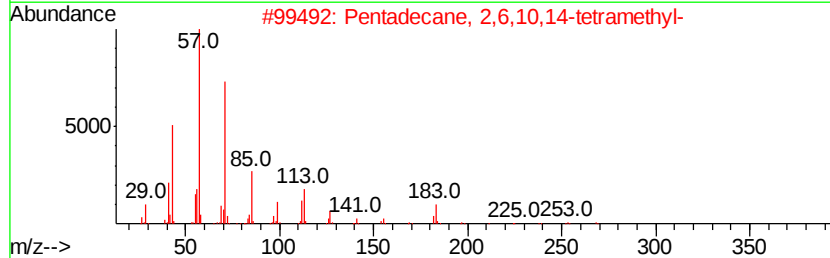
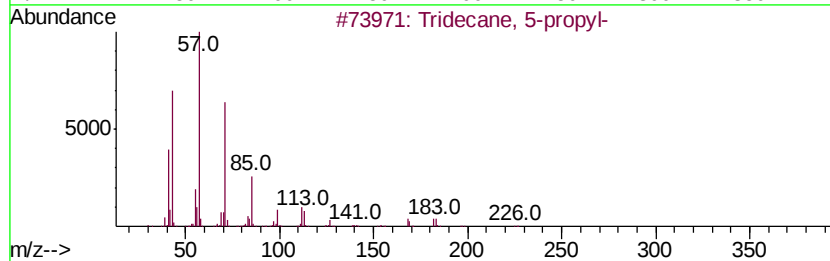
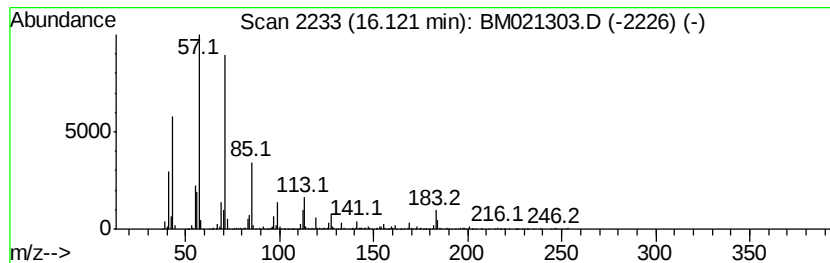
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	15.66 ng/ul	3059500	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	93
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	87
3		Heptacosane	380	C27H56	000593-49-7	80
4		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	74
5		Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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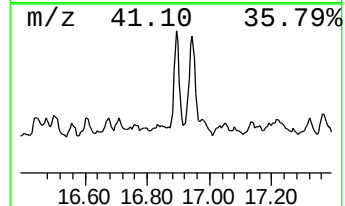
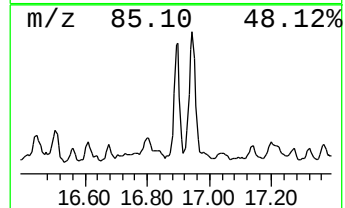
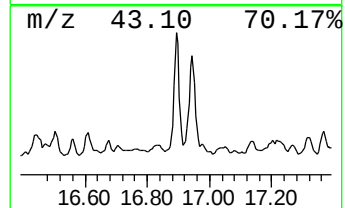
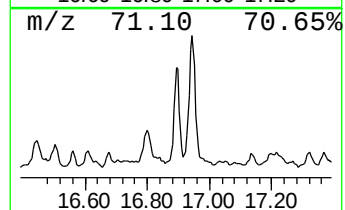
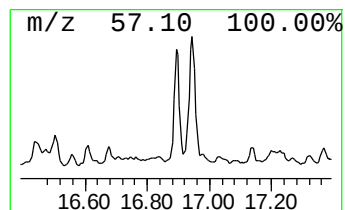
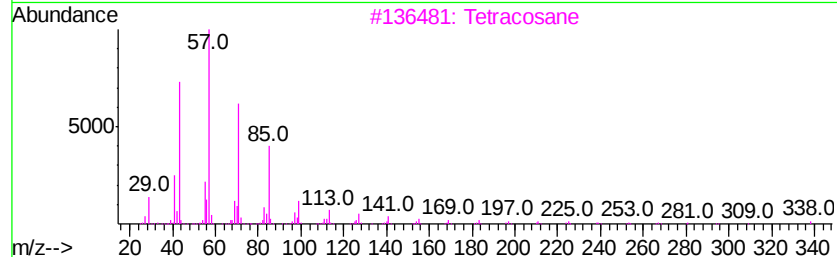
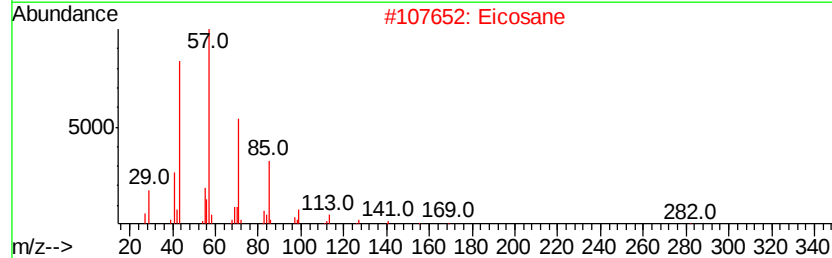
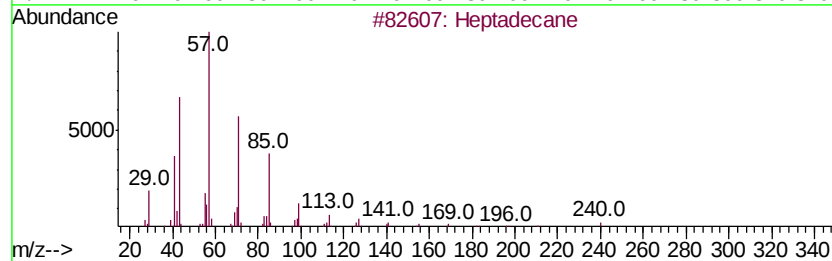
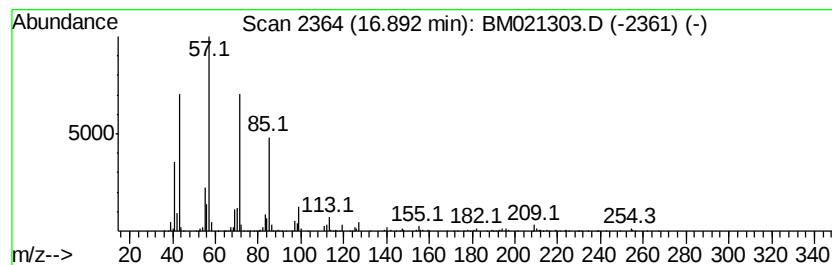
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.89	2.99 ng/ul	585157	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	91
2			Eicosane	282	C20H42	000112-95-8	87
3			Tetracosane	338	C24H50	000646-31-1	87
4			Hexadecane	226	C16H34	000544-76-3	86
5			Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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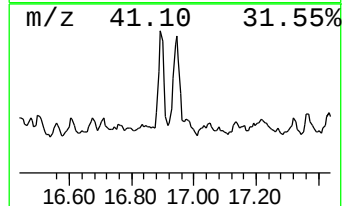
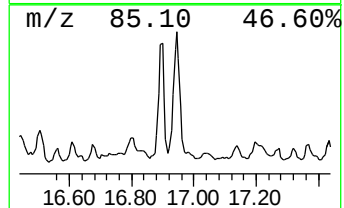
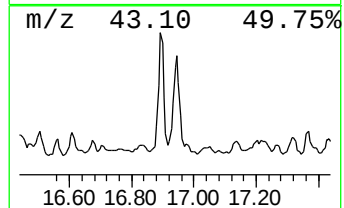
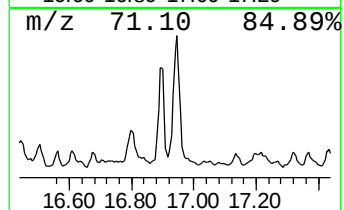
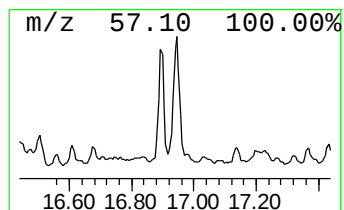
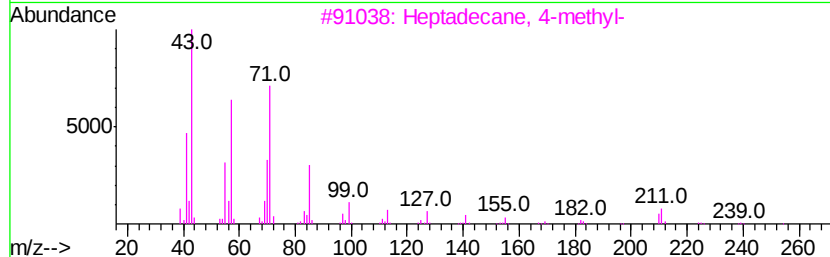
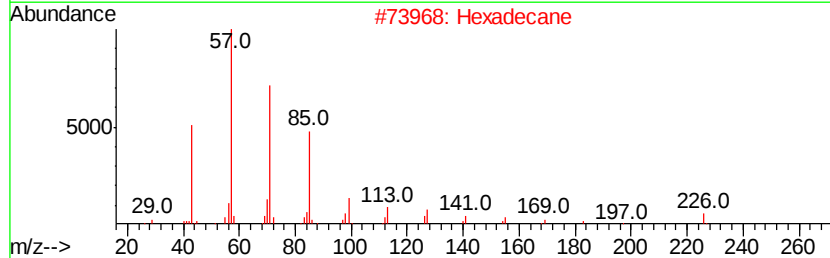
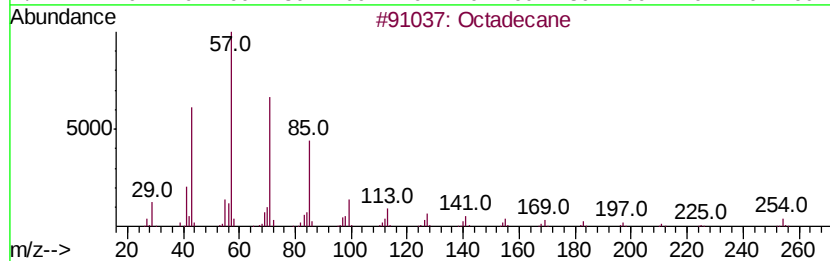
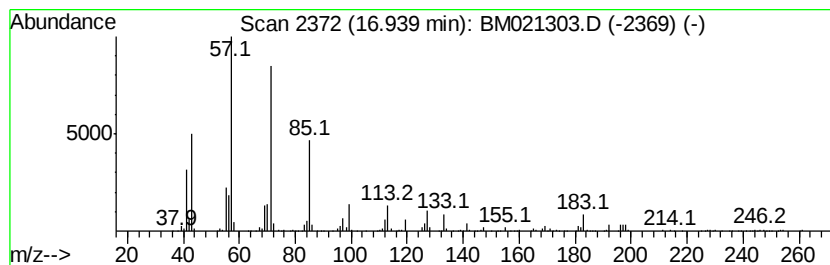
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.94	4.32 ng/ul	844056	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	83
2		Hexadecane	226	C16H34	000544-76-3	80
3		Heptadecane, 4-methyl-	254	C18H38	026429-11-8	76
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	74
5		Hexadecane	226	C16H34	000544-76-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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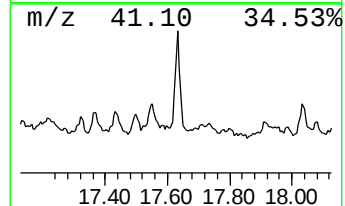
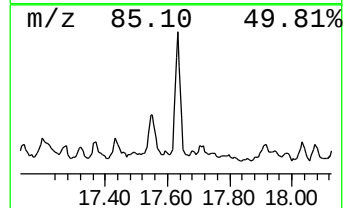
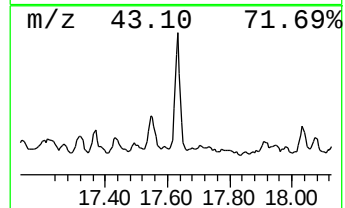
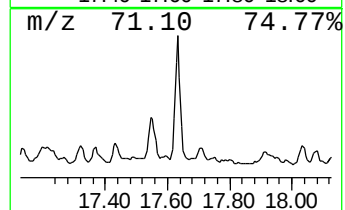
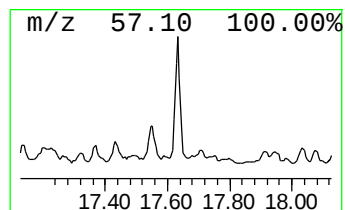
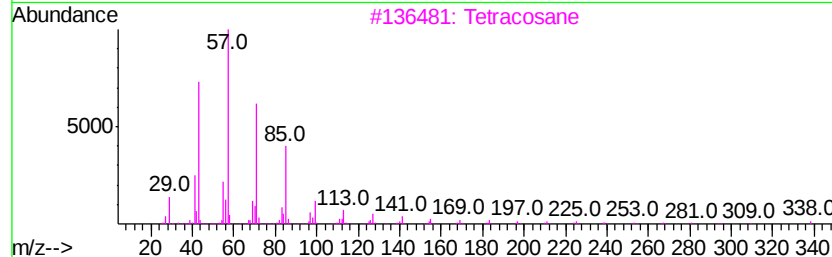
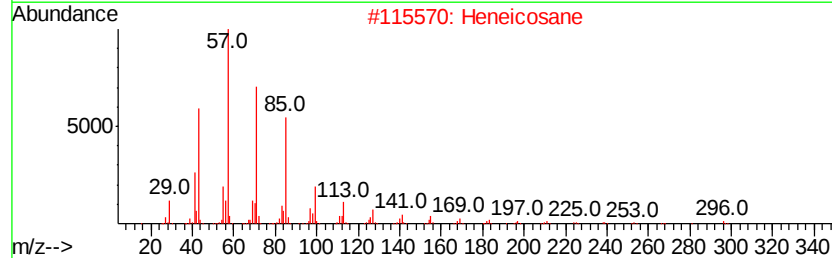
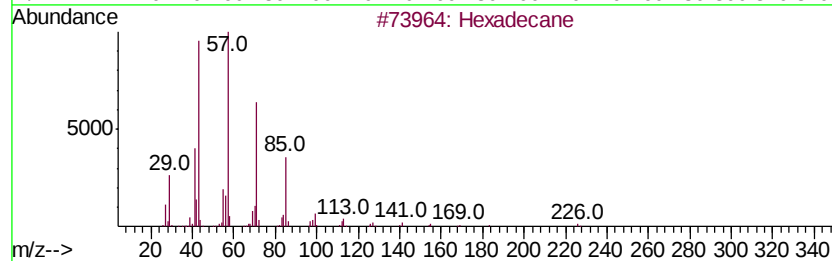
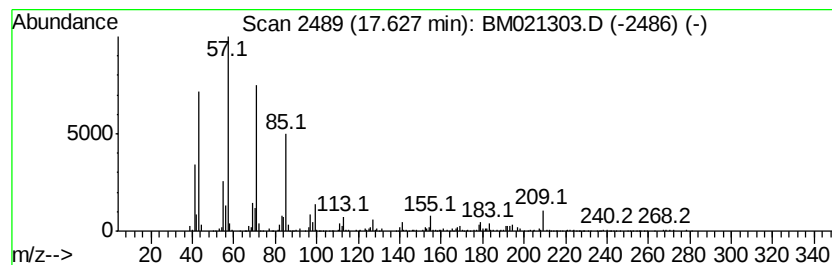
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.63	4.15 ng/ul	811798	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Heneicosane	296	C21H44	000629-94-7	87
3		Tetracosane	338	C24H50	000646-31-1	87
4		Decane, 5-propyl-	184	C13H28	017312-62-8	83
5		Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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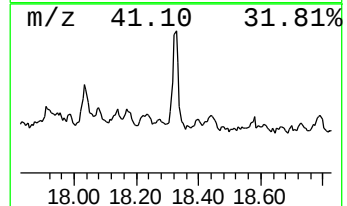
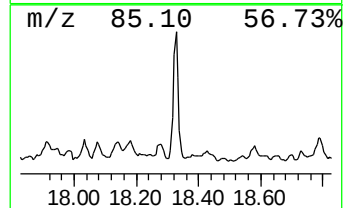
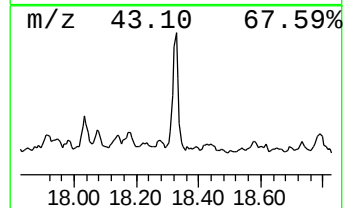
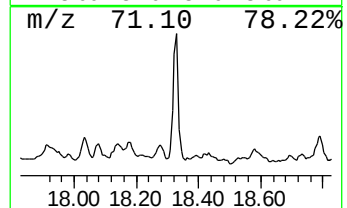
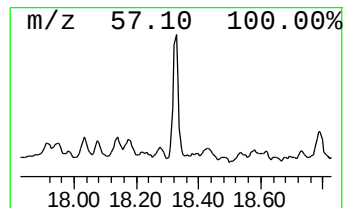
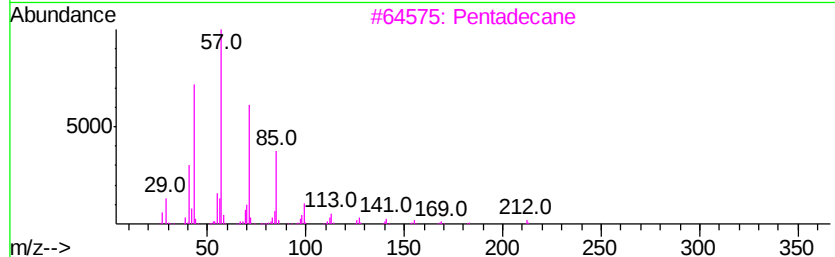
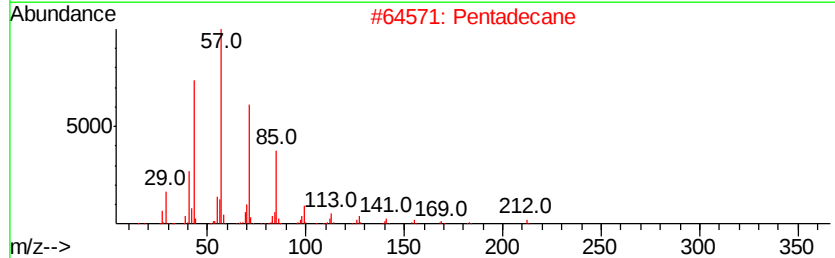
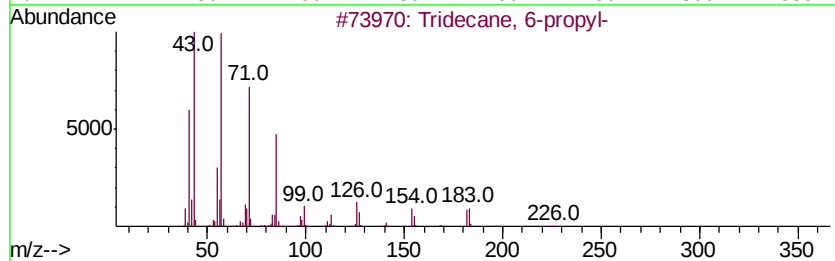
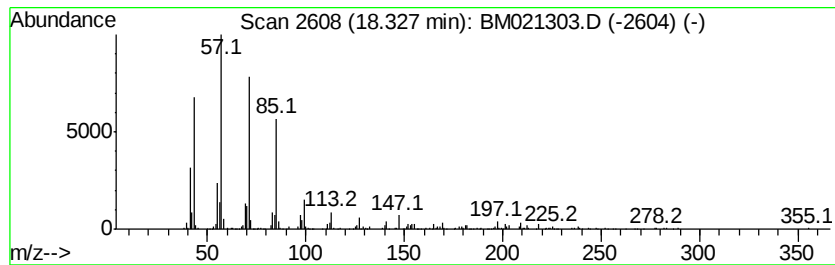
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	3.47 ng/ul	678265	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 6-propyl-	226	C16H34	055045-10-8	91
2			Pentadecane	212	C15H32	000629-62-9	90
3			Pentadecane	212	C15H32	000629-62-9	87
4			Nonadecane	268	C19H40	000629-92-5	87
5			Tridecane, 7-propyl-	226	C16H34	055045-09-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5BA3DL

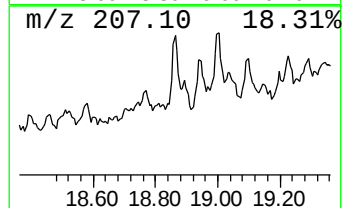
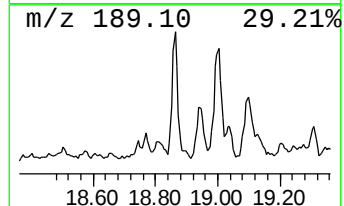
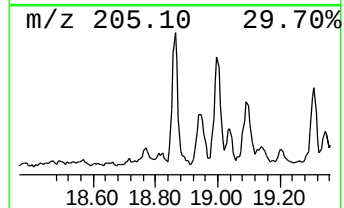
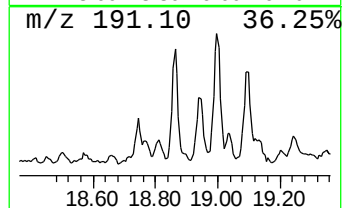
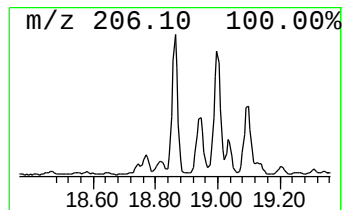
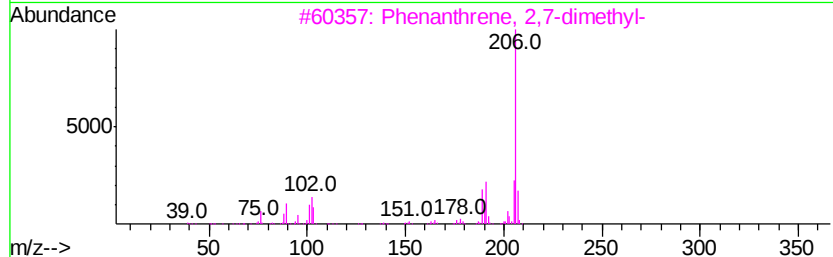
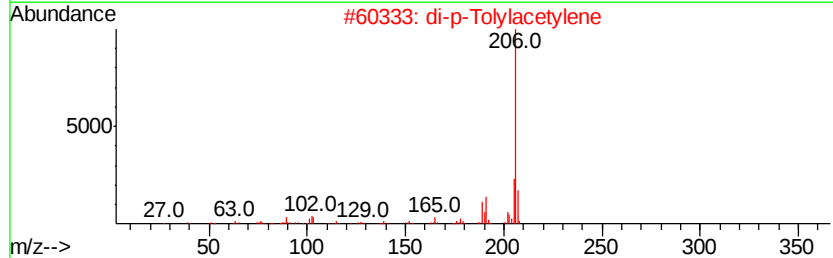
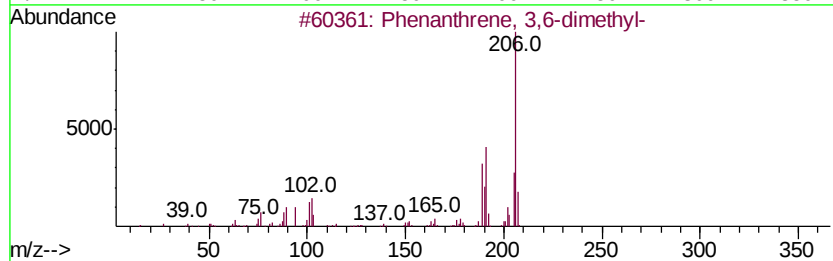
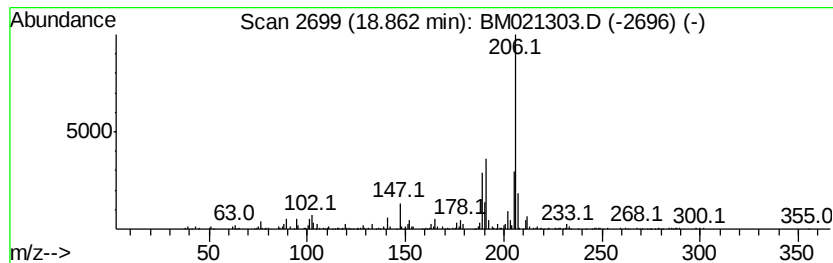
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenanthrene, 3,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	2.35 ng/ul	458438	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	96
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	90
4			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
5			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5BA3DL

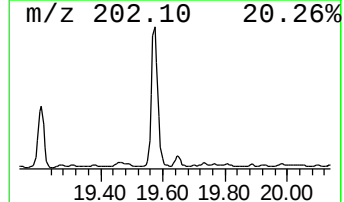
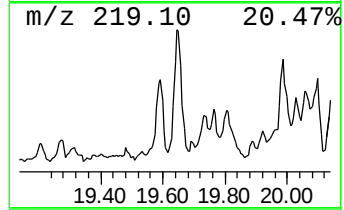
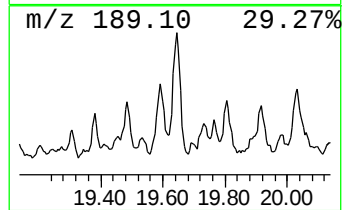
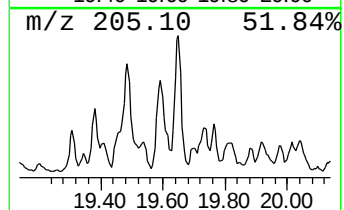
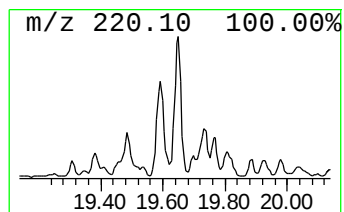
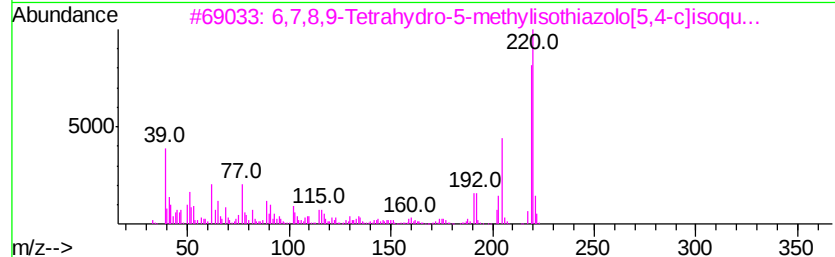
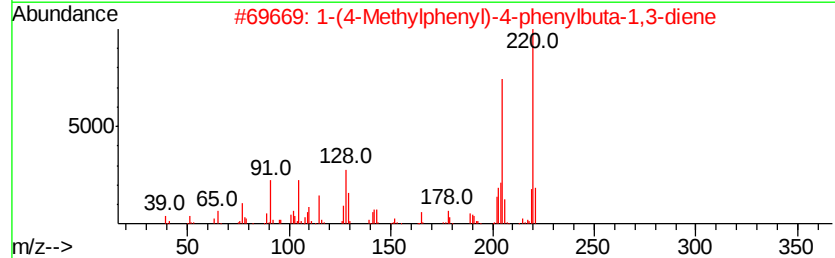
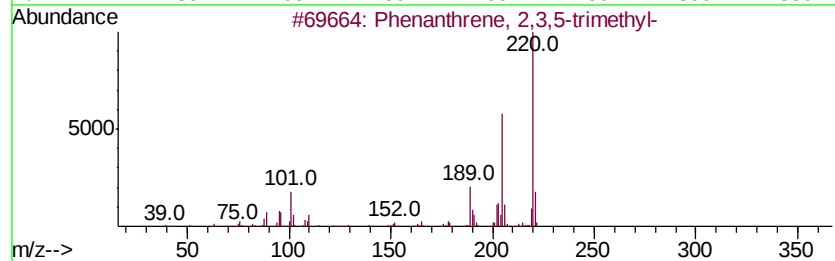
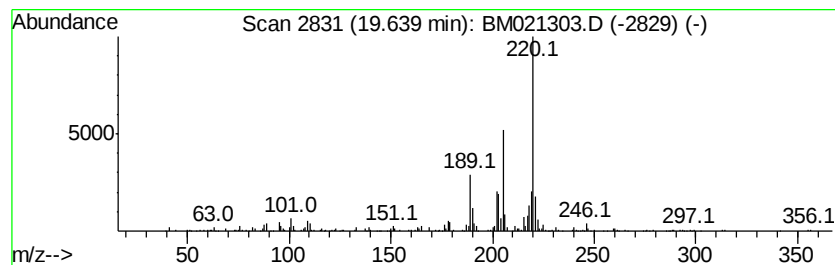
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.64	3.81 ng/ul	575035	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	81
2		1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	74
3		6,7,8,9-Tetrahydro-5-methylisoth...	220	C11H12N2OS	339156-87-5	58
4		Pyrrolo[1,2-a]thieno[2,3-d]pyrim...	220	C11H12N2OS	056929-61-4	58
5		1-Penten-3-one, 4-methyl-1-[2,6,...	220	C15H24O	1000132-20-9	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070119\
Data File : BM021303.D
Acq On : 01 Jul 2019 11:11
Operator : HP/JU
Sample : K3590-04DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
C5BA3DL

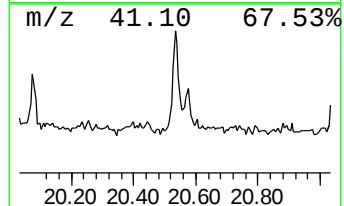
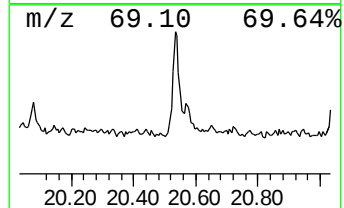
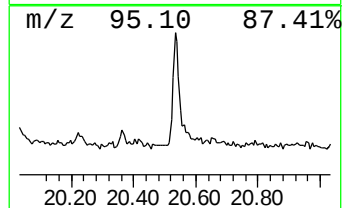
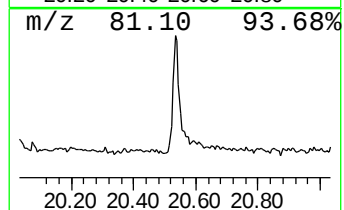
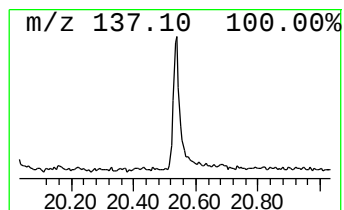
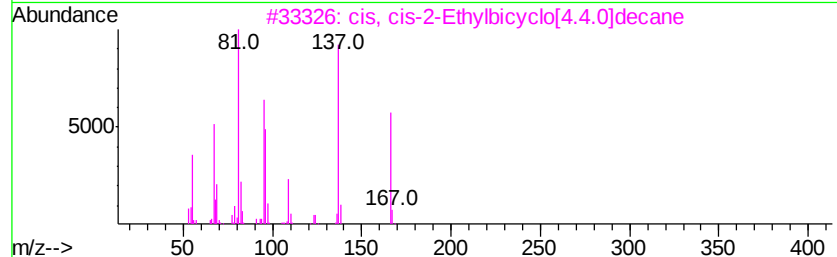
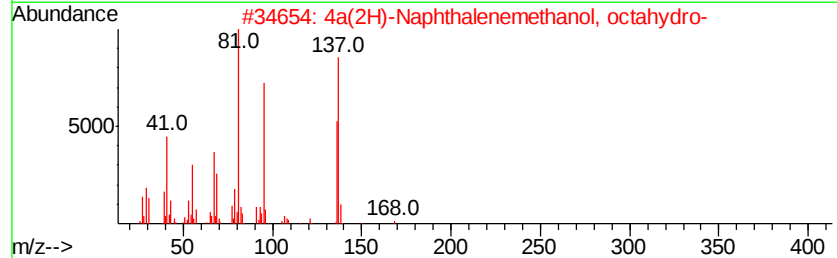
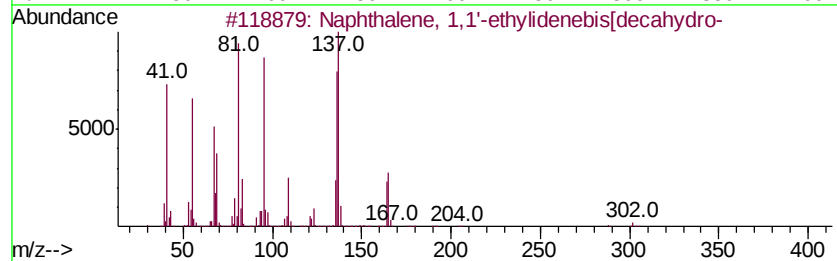
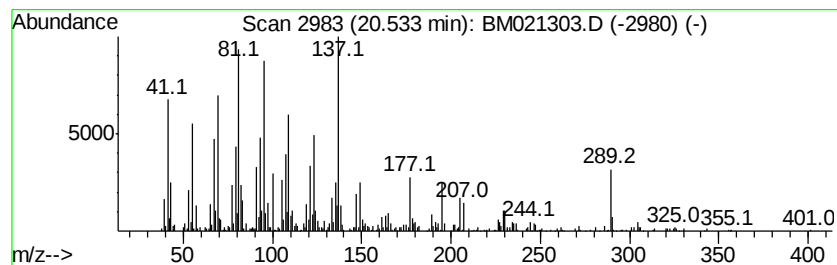
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 Naphthalene, 1,1'-ethyliden... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.53	3.36 ng/ul	507921	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,1'-ethylidenebis[...	302	C22H38	054934-70-2	78
2		4a(2H)-Naphthalenemethanol, octa...	168	C11H20O	099992-19-5	55
3		cis, cis-2-Ethylbicyclo[4.4.0]de...	166	C12H22	066660-40-0	47
4		2,5-Octadiene, 3,4,5,6-tetramethyl-	166	C12H22	247797-85-9	46
5		Trans, trans-2-ethylbicyclo[4.4....	166	C12H22	066660-37-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM070119\
 Data File : BM021303.D
 Acq On : 01 Jul 2019 11:11
 Operator : HP/JU
 Sample : K3590-04DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5BA3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.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
unknown-01	3.01	6.9	ng/ul	451109	1	7.76	1311190	20.0
Benzene, 2,4-diis...	12.77	3.4	ng/ul	541203	3	14.40	3150820	20.0
1,4-Benzenedicarb...	13.46	2.0	ng/ul	318861	3	14.40	3150820	20.0
(DEL) Alkane: Str...	15.24	3.2	ng/ul	499838	3	14.40	3150820	20.0
(DEL) Alkane: Str...	15.64	5.5	ng/ul	866975	3	14.40	3150820	20.0
(DEL) Alkane: Str...	16.12	15.7	ng/ul	3059500	4	17.15	3907930	20.0
(DEL) Alkane: Str...	16.89	3.0	ng/ul	585157	4	17.15	3907930	20.0
(DEL) Alkane: Str...	16.94	4.3	ng/ul	844056	4	17.15	3907930	20.0
(DEL) Alkane: Str...	17.63	4.2	ng/ul	811798	4	17.15	3907930	20.0
(DEL) Alkane: Str...	18.33	3.5	ng/ul	678265	4	17.15	3907930	20.0
Phenanthrene, 3,6...	18.86	2.4	ng/ul	458438	4	17.15	3907930	20.0
Phenanthrene, 2,3...	19.64	3.8	ng/ul	575035	5	21.33	3021530	20.0
Naphthalene, 1,1'...	20.53	3.4	ng/ul	507921	5	21.33	3021530	20.0