

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
 Data File : BP000142.D
 Acq On : 09 Sep 2019 19:13
 Operator : HP/JU
 Sample : K4633-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F2D08

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_P\METHODS\SOM-EPA-BP090519MA.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	2.170	6	14	18	rBV	12143078	19326056	100.00%	12.860%
2	2.611	85	89	97	rVB	81046	113800	0.59%	0.076%
3	3.970	316	320	323	rVV	119904	156521	0.81%	0.104%
4	6.505	747	751	758	rBV2	2529532	2482935	12.85%	1.652%
5	6.569	758	762	766	rBV	2021083	1789254	9.26%	1.191%
6	6.658	773	777	781	rBV	2892937	2312456	11.97%	1.539%
7	6.893	813	817	820	rBV	2843754	2194541	11.36%	1.460%
8	7.258	875	879	882	rBV	3242287	2519460	13.04%	1.676%
9	7.446	907	911	914	rBV	2582901	2017081	10.44%	1.342%
10	7.775	963	967	970	rBV	1952253	1575868	8.15%	1.049%
11	8.016	1004	1008	1011	rBV	2968265	2662146	13.77%	1.771%
12	8.175	1032	1035	1038	rBV	3586222	3033327	15.70%	2.018%
13	8.199	1038	1039	1041	rVV	238433	131152	0.68%	0.087%
14	8.228	1041	1044	1048	rVB	2584979	2252738	11.66%	1.499%
15	9.634	1279	1283	1285	rVV	4104537	3706692	19.18%	2.466%
16	9.652	1285	1286	1290	rVB	1560525	1097684	5.68%	0.730%
17	9.787	1305	1309	1319	rVB	5225084	4742353	24.54%	3.156%
18	9.946	1332	1336	1339	rBV	4664551	3600204	18.63%	2.396%
19	9.975	1339	1341	1344	rVB	544816	413493	2.14%	0.275%
20	10.028	1346	1350	1360	rVV	1604397	1309968	6.78%	0.872%
21	10.146	1367	1370	1374	rVB2	420148	344831	1.78%	0.229%
22	10.469	1421	1425	1428	rBV	6789147	5458655	28.25%	3.632%
23	10.499	1428	1430	1432	rVV	220139	185260	0.96%	0.123%
24	10.522	1432	1434	1438	rVV	1236071	1074759	5.56%	0.715%
25	10.599	1443	1447	1451	rVV2	98011	136807	0.71%	0.091%
26	10.863	1486	1492	1495	rVV3	111444	128155	0.66%	0.085%
27	11.246	1554	1557	1563	rVB	167261	151323	0.78%	0.101%
28	11.340	1571	1573	1579	rVB	120995	108418	0.56%	0.072%
29	11.446	1588	1591	1594	rVV	4063496	4118523	21.31%	2.741%
30	11.469	1594	1595	1598	rVV	2526461	1975334	10.22%	1.314%
31	11.504	1598	1601	1607	rVV	6265319	5978279	30.93%	3.978%
32	11.675	1625	1630	1634	rBV2	407460	374144	1.94%	0.249%
33	11.957	1675	1678	1680	rBV	215462	179284	0.93%	0.119%
34	11.999	1680	1685	1688	rBV2	1040797	1092907	5.66%	0.727%

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35	12.034	1688	1691	1694	rVB2	110171	119676	0.62%	0.080%
36	12.075	1694	1698	1700	rBV	389540	414824	2.15%	0.276%
37	12.257	1726	1729	1731	rBV	175985	180927	0.94%	0.120%
38	12.275	1731	1732	1736	rVB	281070	203325	1.05%	0.135%
39	12.516	1769	1773	1777	rBV2	123503	159321	0.82%	0.106%
40	12.599	1785	1787	1793	rVB2	161018	172469	0.89%	0.115%
41	12.681	1793	1801	1805	rBV	4872647	4580398	23.70%	3.048%
42	12.740	1805	1811	1815	rVB2	69091	105327	0.54%	0.070%
43	12.799	1818	1821	1823	rBV	265235	269722	1.40%	0.179%
44	12.822	1823	1825	1834	rVB2	162413	269667	1.40%	0.179%
45	12.899	1834	1838	1840	rBV	7052609	7590544	39.28%	5.051%
46	12.963	1846	1849	1856	rVB	175526	209861	1.09%	0.140%
47	13.040	1858	1862	1864	rBV	209269	198056	1.02%	0.132%
48	13.069	1864	1867	1873	rVB2	305488	311918	1.61%	0.208%
49	13.140	1873	1879	1882	rBV3	236848	399714	2.07%	0.266%
50	13.228	1891	1894	1895	rVV2	217182	216791	1.12%	0.144%
51	13.251	1895	1898	1901	rVB	512347	524413	2.71%	0.349%
52	13.322	1906	1910	1912	rBV	315652	340275	1.76%	0.226%
53	13.357	1912	1916	1918	rVV	446078	412345	2.13%	0.274%
54	13.381	1918	1920	1924	rVB	174400	166784	0.86%	0.111%
55	13.451	1929	1932	1935	rBV	239828	214247	1.11%	0.143%
56	13.481	1935	1937	1941	rBV	201508	190437	0.99%	0.127%
57	13.557	1947	1950	1956	rBV5	73571	131074	0.68%	0.087%
58	13.663	1960	1968	1969	rBV6	136467	258585	1.34%	0.172%
59	13.687	1969	1972	1975	rVB2	174175	197764	1.02%	0.132%
60	13.763	1982	1985	1991	rVB	376417	450933	2.33%	0.300%
61	13.881	2000	2005	2008	rBV2	598347	756380	3.91%	0.503%
62	13.916	2008	2011	2012	rBV	246345	206653	1.07%	0.138%
63	13.934	2012	2014	2019	rVB2	276123	324566	1.68%	0.216%
64	13.981	2019	2022	2024	rVB2	343146	336154	1.74%	0.224%
65	14.016	2024	2028	2033	rBV2	993642	1513053	7.83%	1.007%
66	14.140	2041	2049	2052	rBV2	4937579	7730026	40.00%	5.144%
67	14.169	2052	2054	2060	rVB	3187714	2896678	14.99%	1.927%
68	14.251	2062	2068	2070	rBV2	435943	469673	2.43%	0.313%
69	14.281	2071	2073	2075	rVV2	123013	115503	0.60%	0.077%
70	14.304	2075	2077	2079	rVV2	118288	132574	0.69%	0.088%
71	14.334	2079	2082	2084	rVV	276897	303772	1.57%	0.202%

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72	14.363	2084	2087	2092	rVV2	510340	758472	3.92%	0.505%
73	14.404	2092	2094	2101	rVB2	152030	175372	0.91%	0.117%
74	14.469	2102	2105	2108	rVB2	183964	188594	0.98%	0.125%
75	14.504	2108	2111	2112	rBV	174239	156313	0.81%	0.104%
76	14.540	2112	2117	2120	rVV	534584	694014	3.59%	0.462%
77	14.575	2120	2123	2127	rVV2	399695	397056	2.05%	0.264%
78	14.634	2130	2133	2135	rVV2	230410	218611	1.13%	0.145%
79	14.669	2135	2139	2146	rVV5	521059	1014422	5.25%	0.675%
80	14.857	2167	2171	2173	rBV2	194940	257687	1.33%	0.171%
81	15.004	2194	2196	2199	rVB	166576	176167	0.91%	0.117%
82	15.134	2214	2218	2220	rBV3	181724	211637	1.10%	0.141%
83	15.234	2229	2235	2238	rBV	3672919	6494102	33.60%	4.321%
84	15.340	2250	2253	2256	rVV	511912	567856	2.94%	0.378%
85	15.375	2256	2259	2265	rVB2	311311	532370	2.75%	0.354%
86	15.504	2278	2281	2283	rBV3	92286	110726	0.57%	0.074%
87	15.551	2284	2289	2292	rVV	2230877	3257346	16.85%	2.167%
88	15.587	2292	2295	2298	rVV	3604936	5125888	26.52%	3.411%
89	15.622	2298	2301	2306	rVV	2538246	3088228	15.98%	2.055%
90	15.687	2306	2312	2315	rVV	2355565	3365046	17.41%	2.239%
91	15.716	2315	2317	2325	rVV2	818808	1013567	5.24%	0.674%
92	15.798	2325	2331	2336	rVB3	317236	572412	2.96%	0.381%
93	16.281	2408	2413	2419	rBV3	334046	557210	2.88%	0.371%
94	16.834	2502	2507	2513	rBV3	195903	524466	2.71%	0.349%
95	17.063	2540	2546	2554	rVB4	324235	819240	4.24%	0.545%
96	17.245	2570	2577	2586	rVB2	1416941	3191700	16.52%	2.124%
97	17.334	2588	2592	2599	rVB3	114417	233825	1.21%	0.156%
98	17.487	2604	2618	2624	rBV3	444924	1880502	9.73%	1.251%
99	17.710	2648	2656	2670	rVB	1042811	2362927	12.23%	1.572%
100	17.951	2691	2697	2704	rVB2	226045	518026	2.68%	0.345%

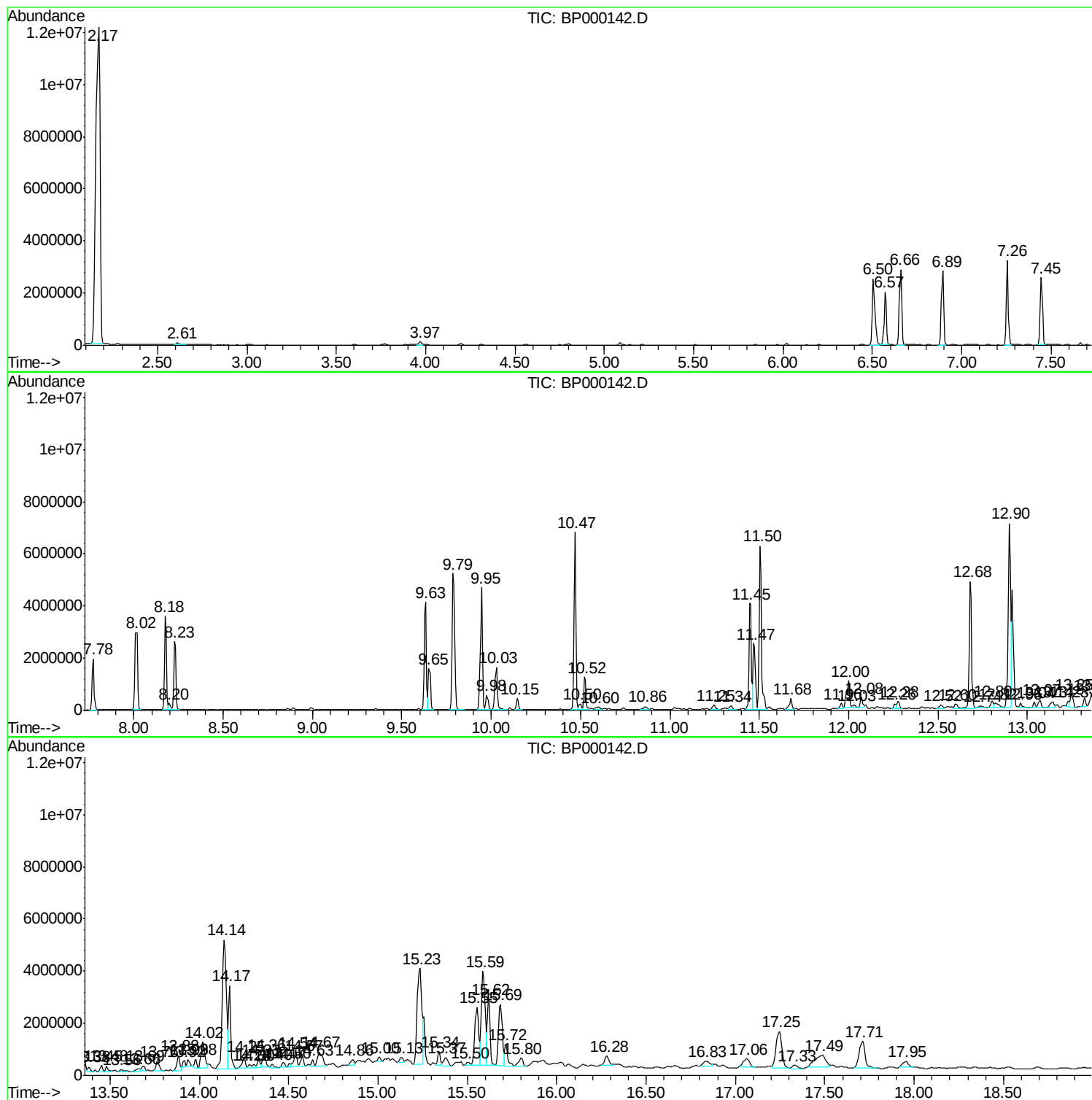
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION

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



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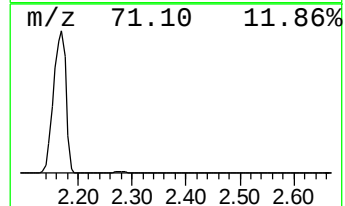
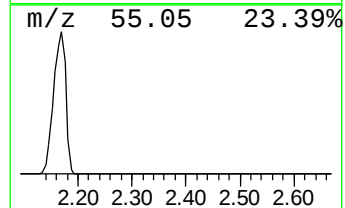
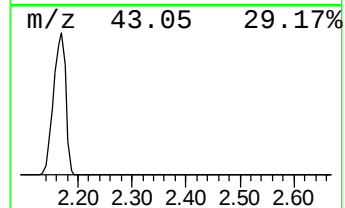
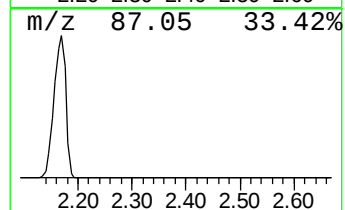
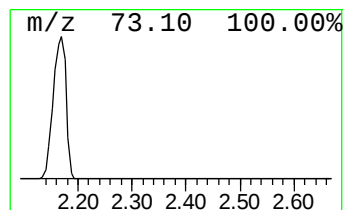
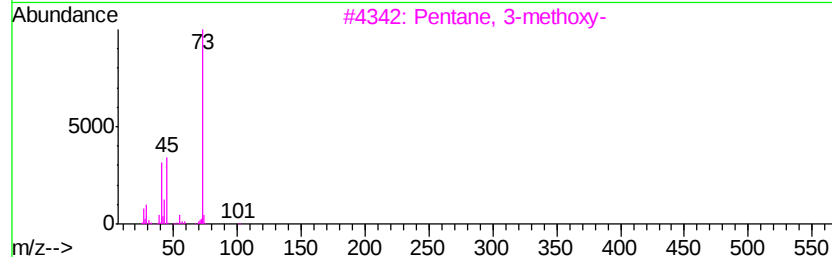
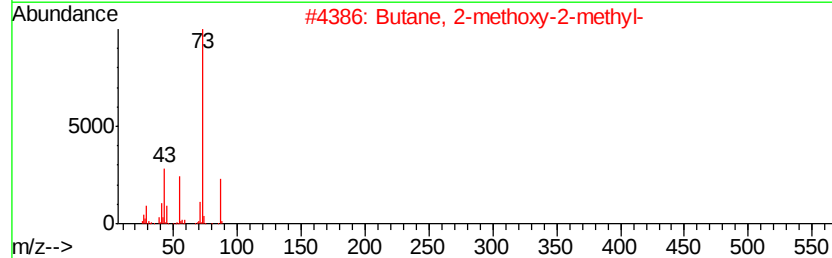
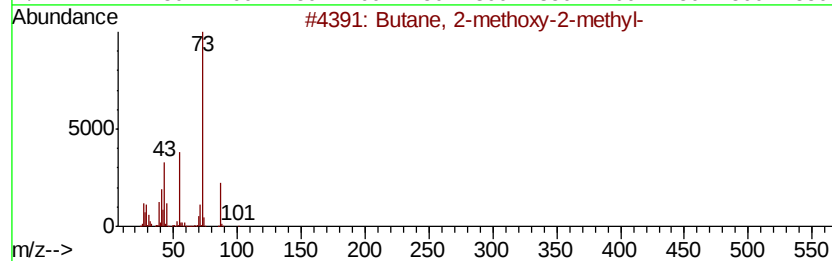
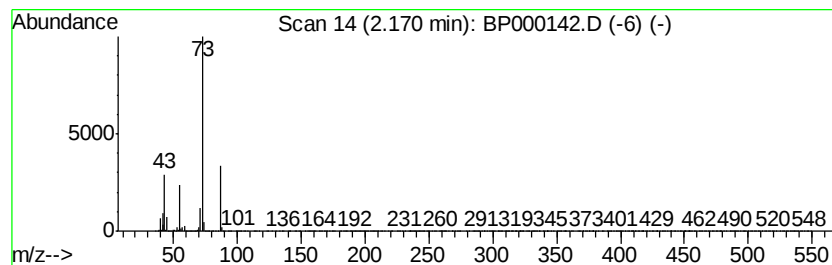
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.17	176.13 ng/ul	19326100	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		1,4-Butanediamine	88	C4H12N2	000110-60-1	9



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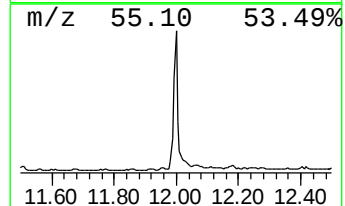
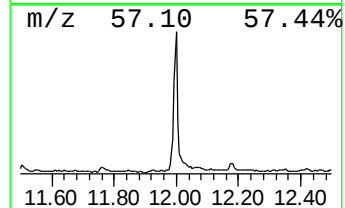
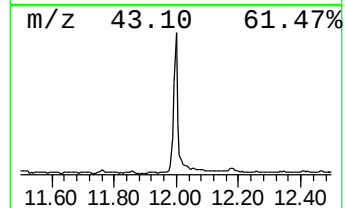
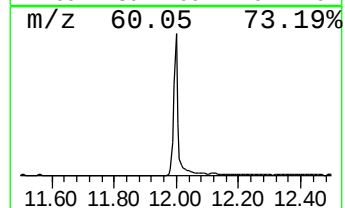
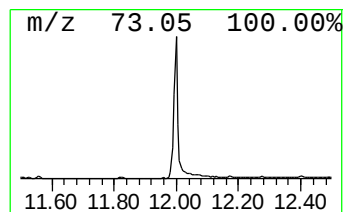
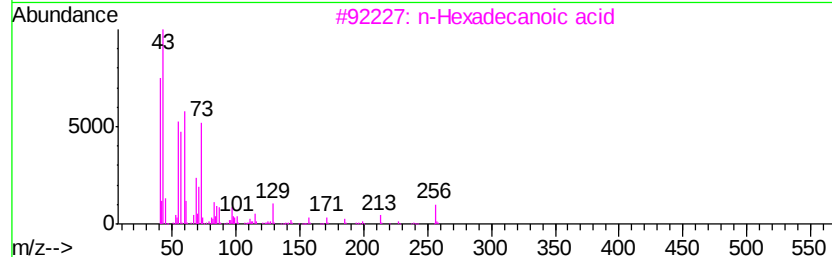
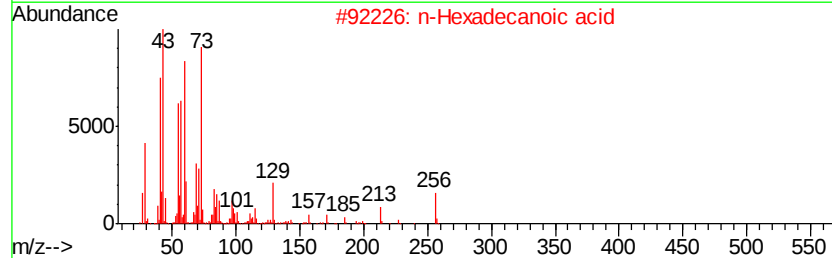
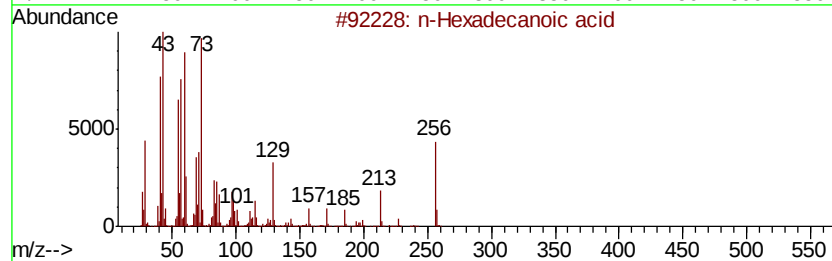
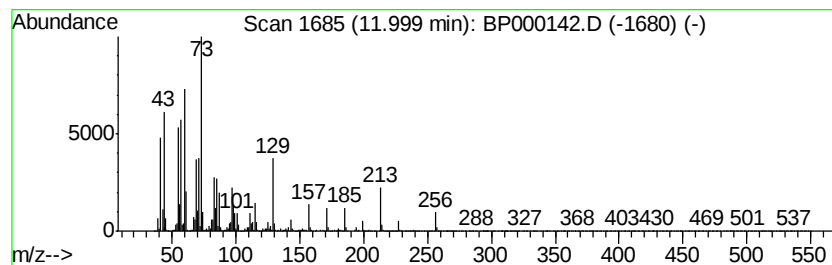
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	5.31 ng/ul	1092910	Phenanthrene-d10	11.45

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96	
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96	
4	Tridecanoic acid	214	C13H26O2	000638-53-9	90	
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	68	



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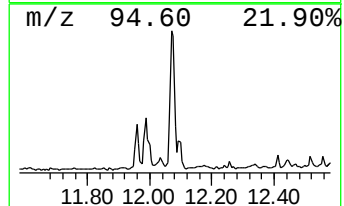
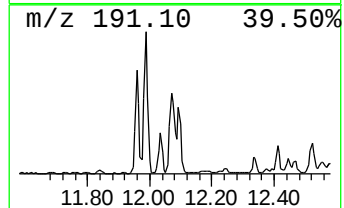
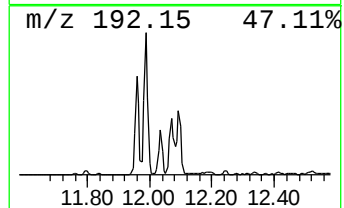
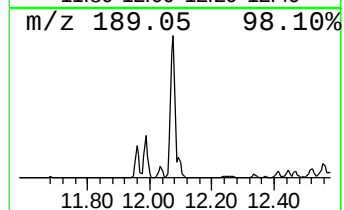
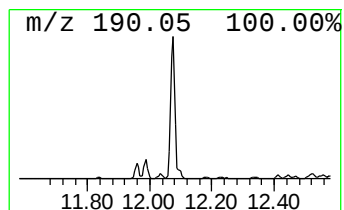
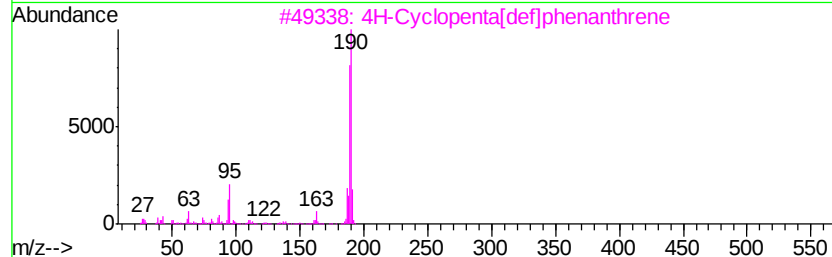
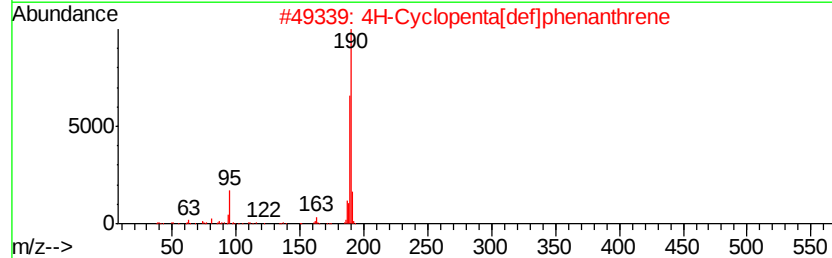
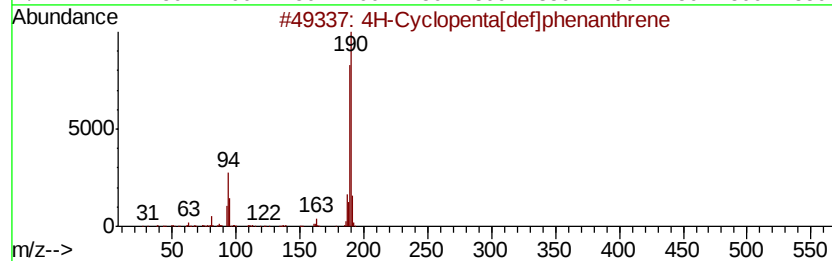
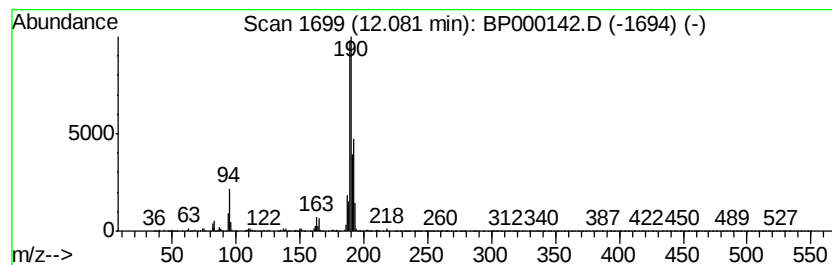
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.08	2.01 ng/ul	414824	Phenanthrene-d10	11.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
5		Methyl diselenide	190	C2H6Se2	007101-31-7	49



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Acq On : 09 Sep 2019 19:13
Operator : HP/JU
Sample : K4633-10
Misc :
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ClientSampleId :
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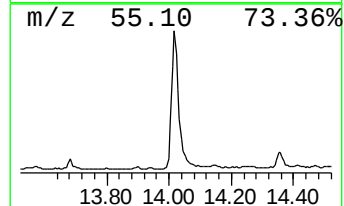
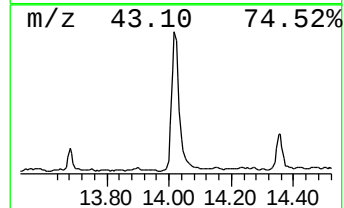
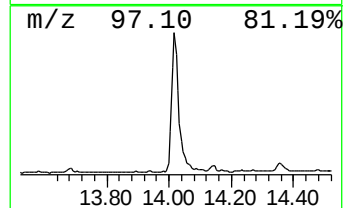
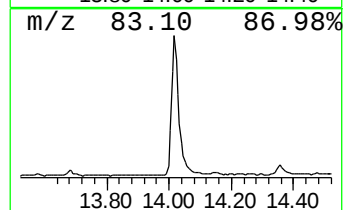
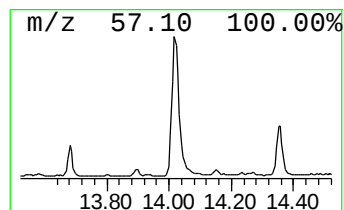
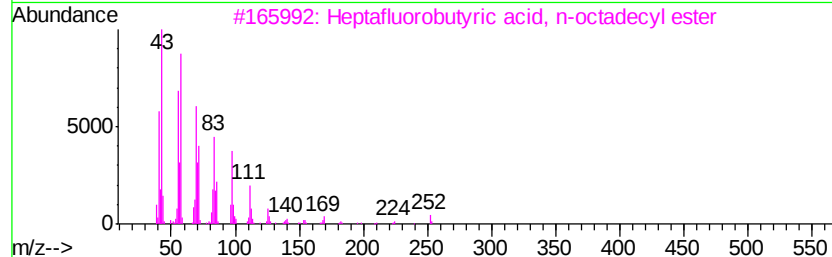
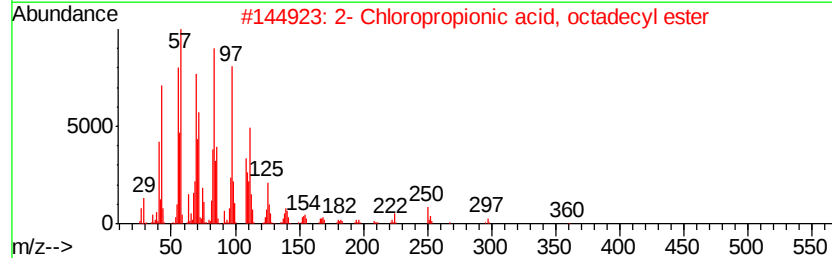
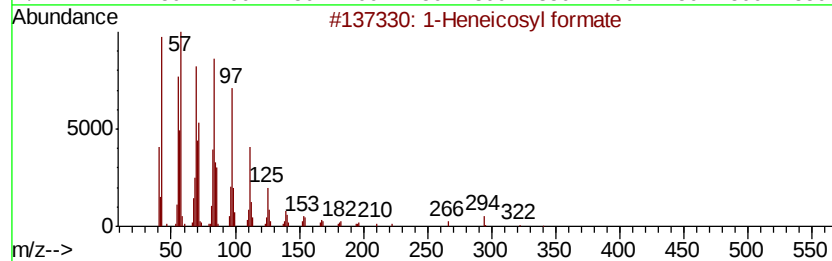
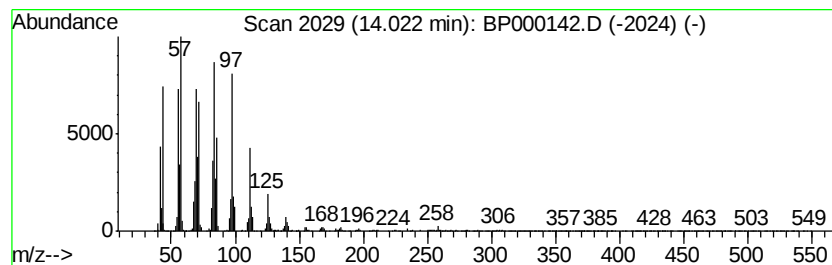
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	3.91 ng/ul	1513050	Chrysene-d12	14.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate		340	C22H44O2	077899-03-7	96
2	2- Chloropropionic acid, octadec...		360	C21H41ClO2	088104-31-8	93
3	Heptafluorobutyric acid, n-octad...		466	C22H37F7O2	000400-57-7	91
4	17-Pentatriacontene		491	C35H70	006971-40-0	91
5	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP090919\
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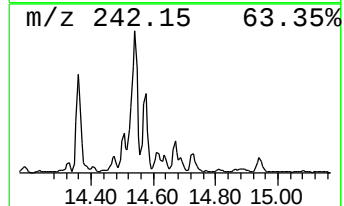
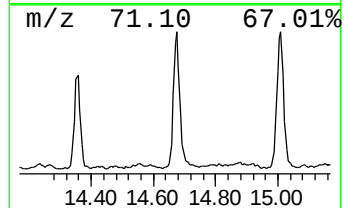
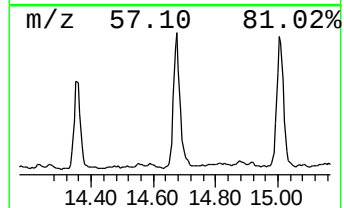
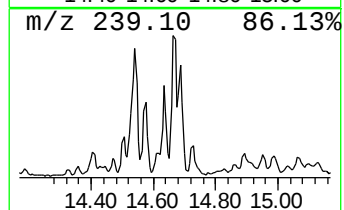
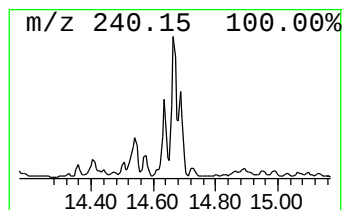
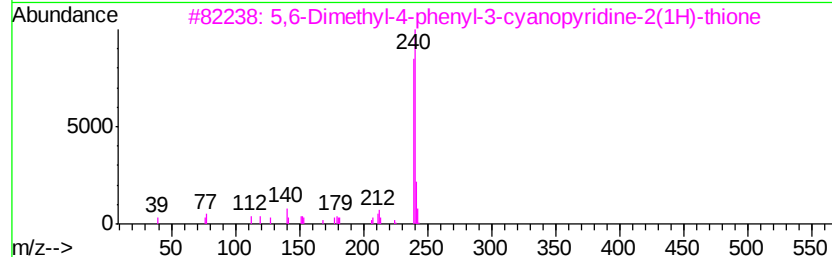
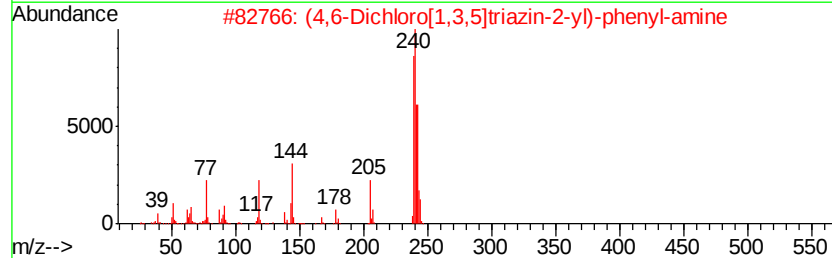
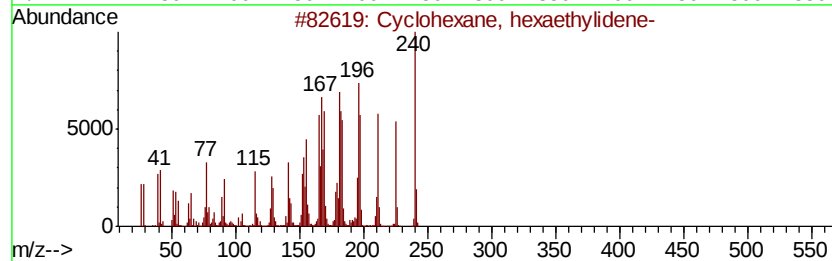
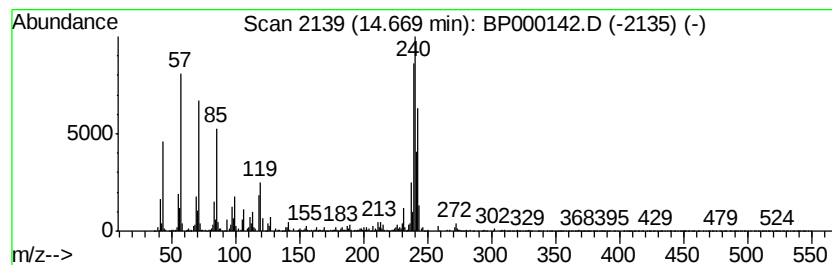
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclohexane, hexaethylidene- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.62 ng/ul	1014420	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, hexaethylidene-	240	C18H24	001482-93-5	64
2		(4,6-Dichloro[1,3,5]triazin-2-yl)-phenyl-amine	240	C9H6Cl2N4	002272-40-4	47
3		5,6-Dimethyl-4-phenyl-3-cyanopyridine	240	C14H12N2S	094639-18-6	25
4		1,3-Dichloro-7-methyl-5,6,7,8-tetrahydrothieno[3,4-c]isoxazole, 6-methyl-	241	C10H9Cl2N3	1000274-39-1	16
5		(3H,6H)Thieno[3,4-c]isoxazole, 6-methyl-	239	C11H10ClNOS	1000115-55-6	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
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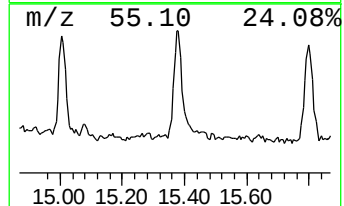
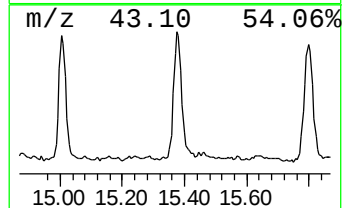
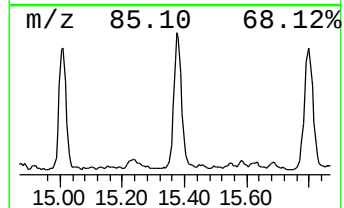
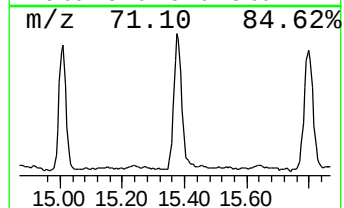
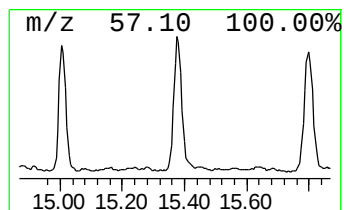
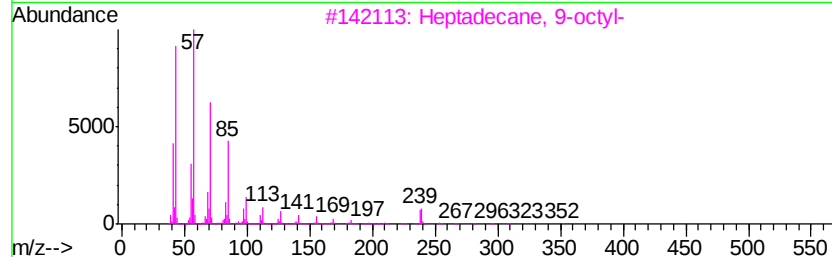
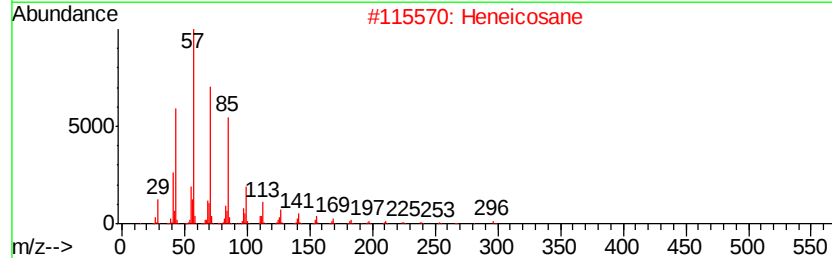
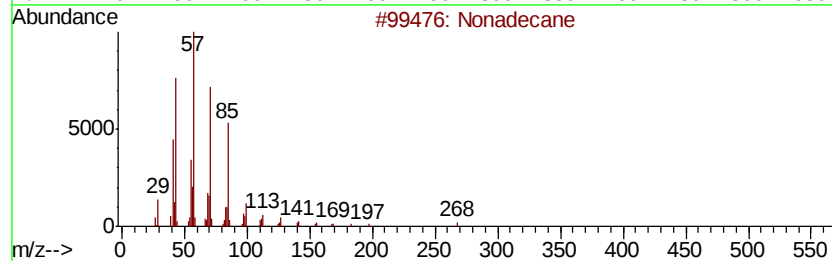
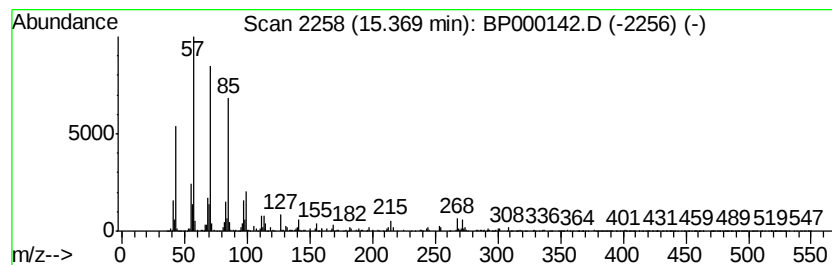
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 Quant Title : SVOA CALIBRATION

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

 Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	3.16 ng/ul	532370	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	86
2		Heneicosane	296	C21H44	000629-94-7	83
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	81
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	81
5		Octadecane	254	C18H38	000593-45-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
Data File : BP000142.D
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Misc :
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ClientSampleId :
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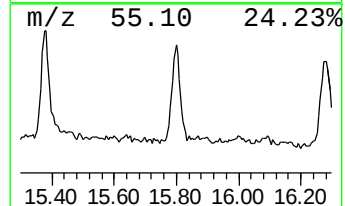
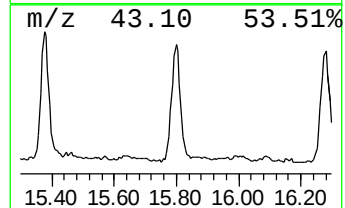
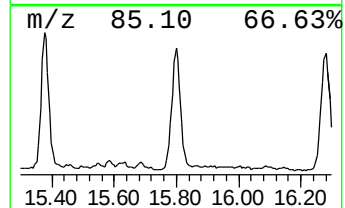
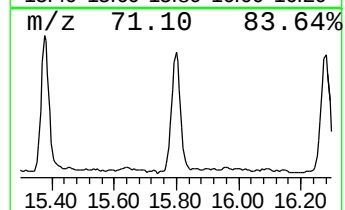
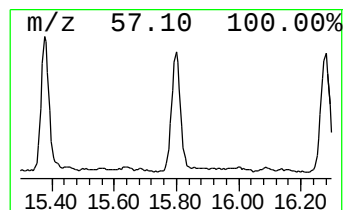
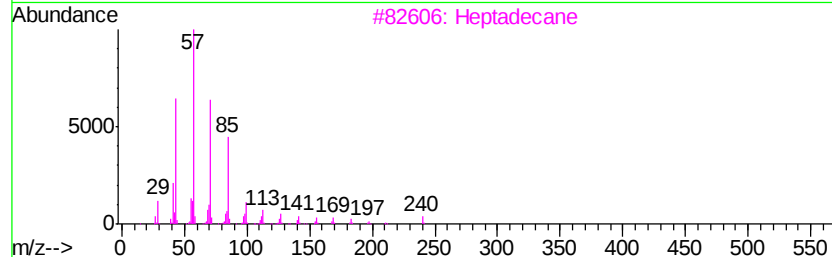
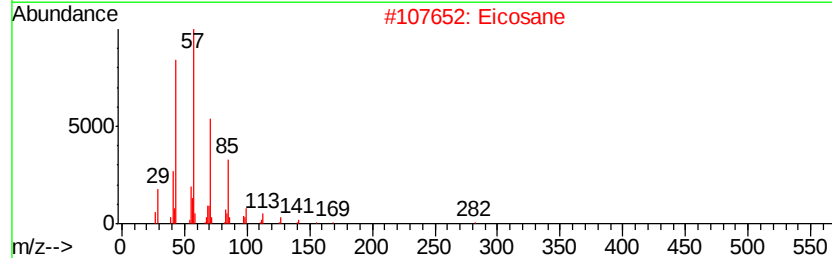
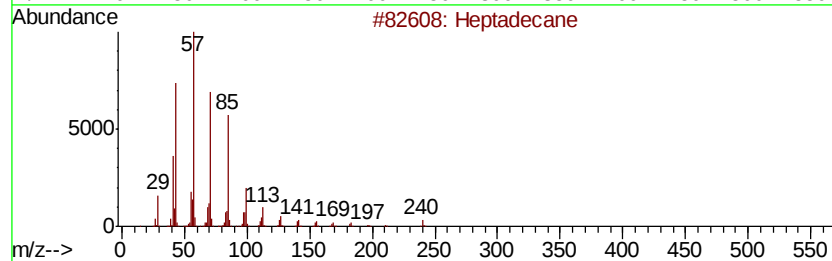
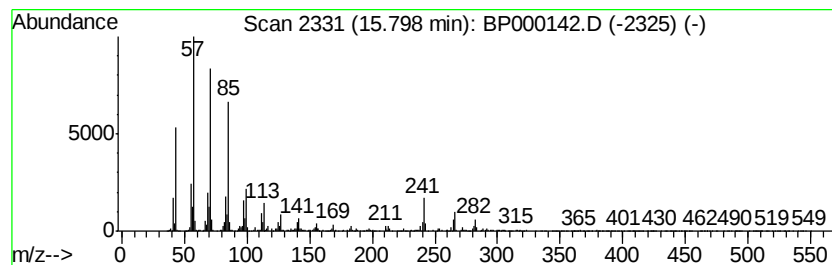
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	3.40 ng/ul	572412	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	98
2		Eicosane	282	C20H42	000112-95-8	96
3		Heptadecane	240	C17H36	000629-78-7	96
4		Eicosane	282	C20H42	000112-95-8	94
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
Data File : BP000142.D
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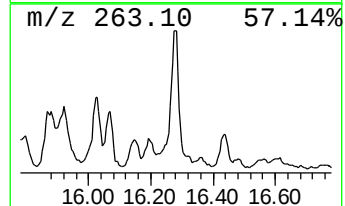
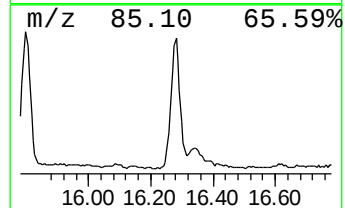
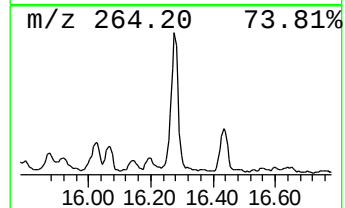
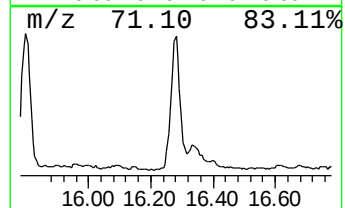
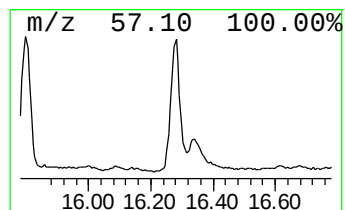
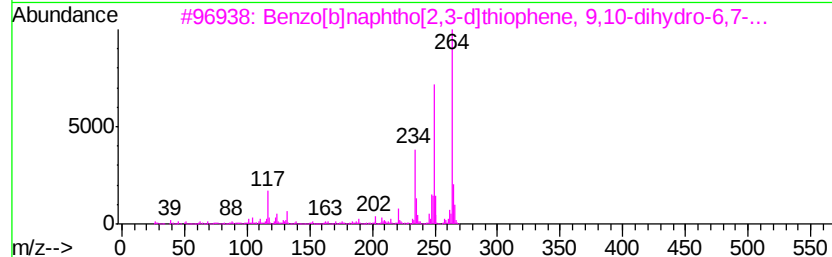
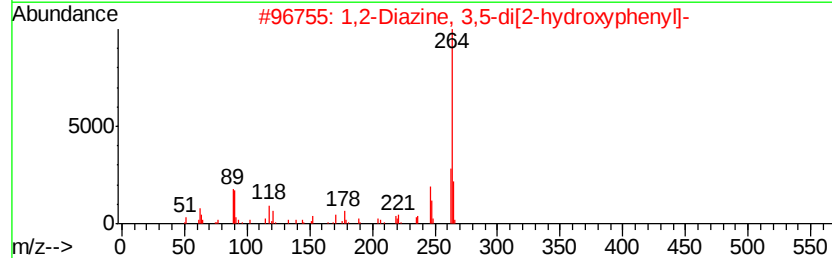
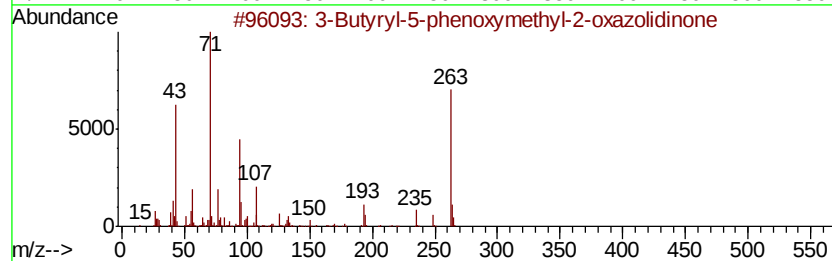
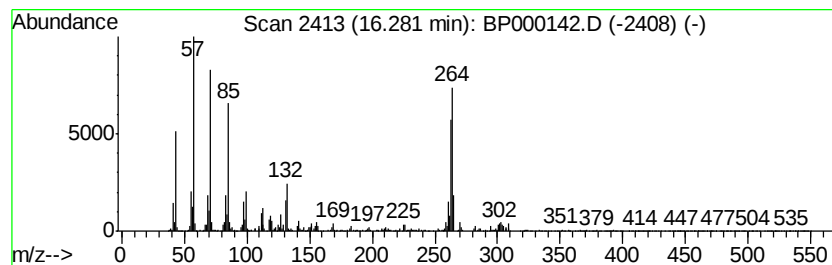
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	3.31 ng/ul	557210	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Butyryl-5-phenoxyethyl-2-oxaz...	263	C14H17N04	014789-96-9	17
2		1,2-Diazine, 3,5-di[2-hydroxyphe...	264	C16H12N2O2	122763-54-6	10
3		Benzo[b]naphtho[2,3-d]thiophene,...	264	C18H16S	024964-02-1	10
4		3,6-Di-m-tolyl-1,2-dihydro-1,2,4...	264	C16H16N4	068614-68-6	10
5		Pyrazolo[1,5-a]pyrimidin-7-amine...	264	C14H12N6	1000276-36-9	10



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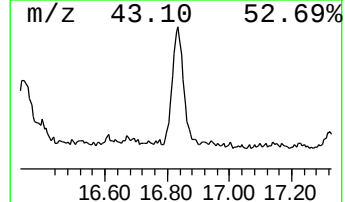
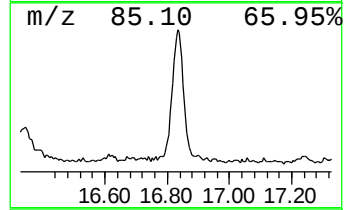
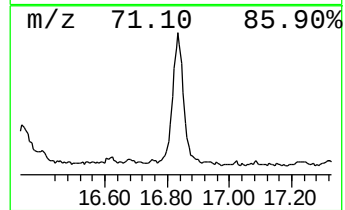
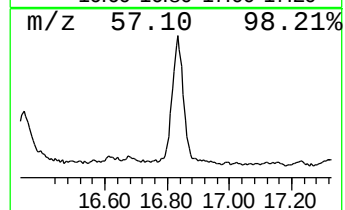
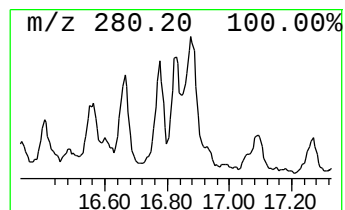
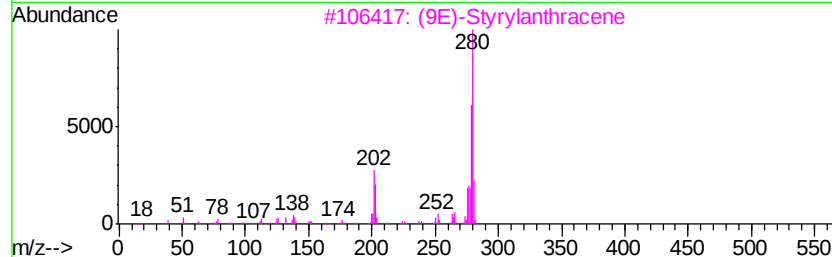
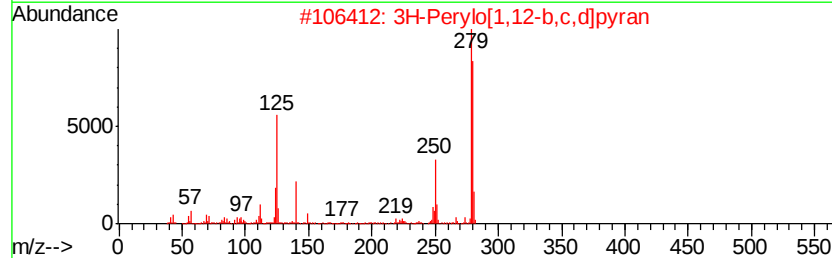
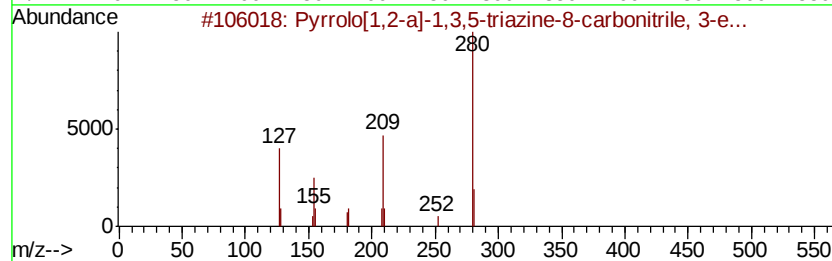
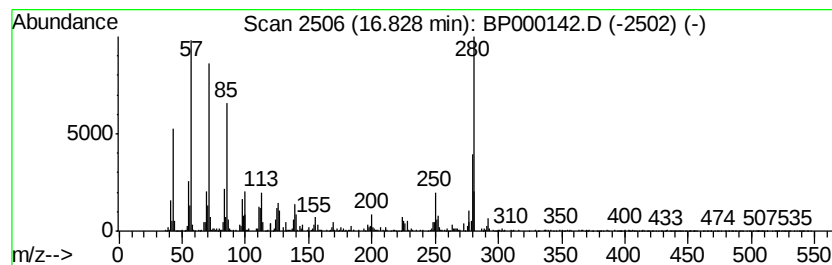
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.83	3.12 ng/ul	524466	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrrolo[1,2-a]-1,3,5-triazine-8-...	280	C15H12N4O2	054450-43-0	32
2		3H-Perylo[1,12-b,c,d]pyran	280	C21H12O	1000281-21-7	30
3		(9E)-Styrylanthracene	280	C22H16	001895-98-3	27
4		Dibenzo(a,c)fluoren-13-one	280	C21H12O	063041-47-4	25
5		Eicosane, 9-octyl-	394	C28H58	013475-77-9	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
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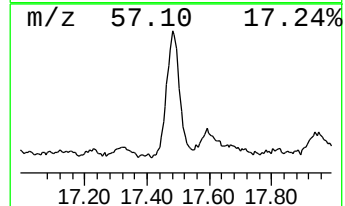
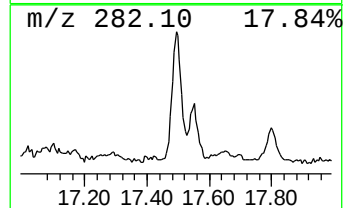
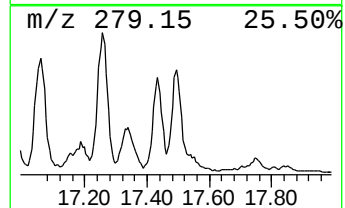
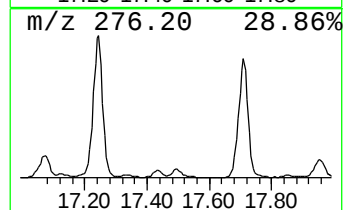
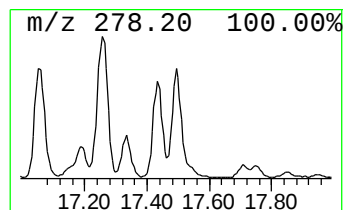
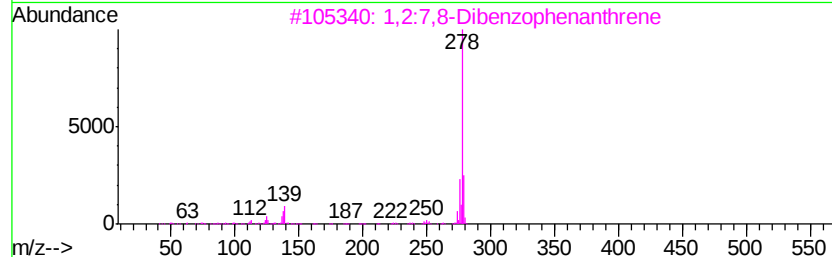
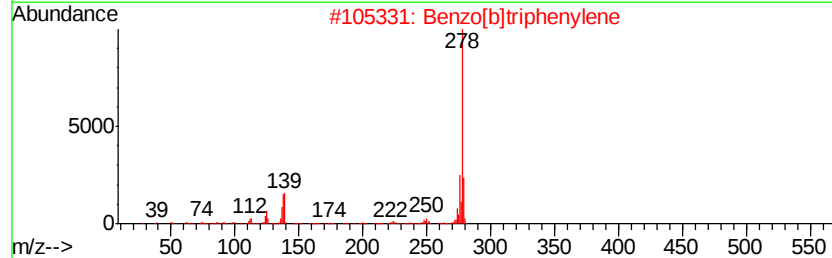
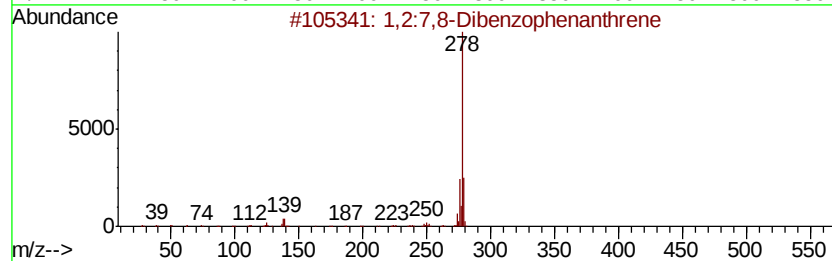
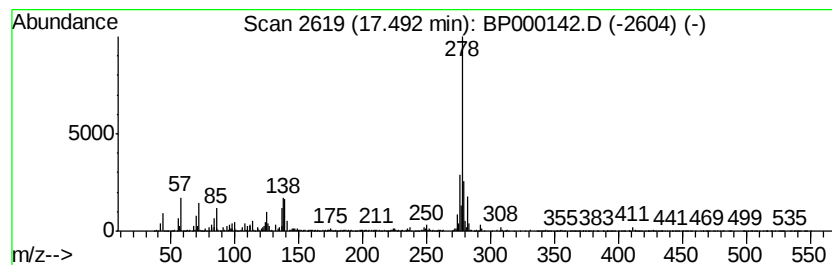
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 1,2:7,8-Dibenzophenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.49	11.18 ng/ul	1880500	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	96
2		Benzo[b]triphenylene	278	C22H14	000215-58-7	93
3		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	92
4		Naphtho[1,2-a]anthracene	278	C22H14	000195-06-2	90
5		Pentaphene	278	C22H14	000222-93-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090919\
Data File : BP000142.D
Acq On : 09 Sep 2019 19:13
Operator : HP/JU
Sample : K4633-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F2D08

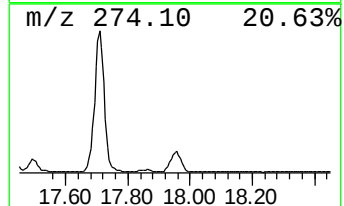
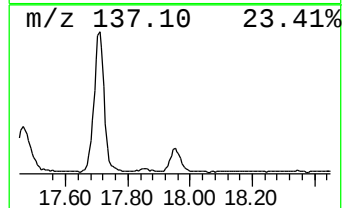
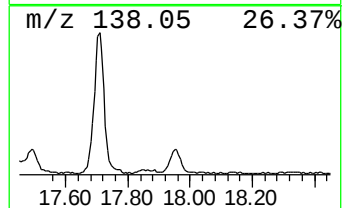
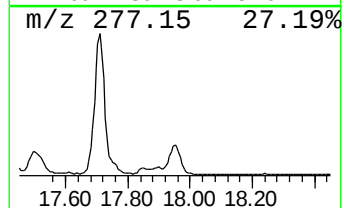
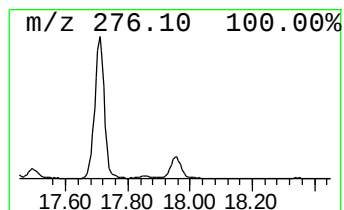
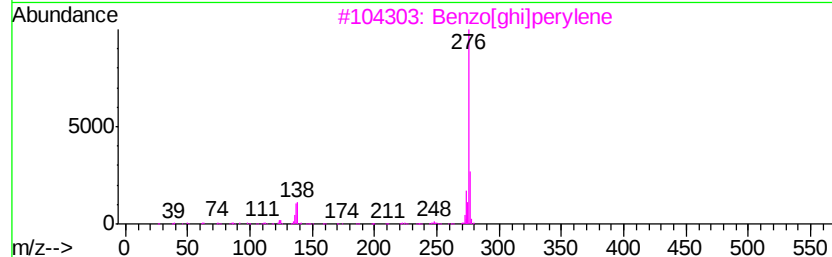
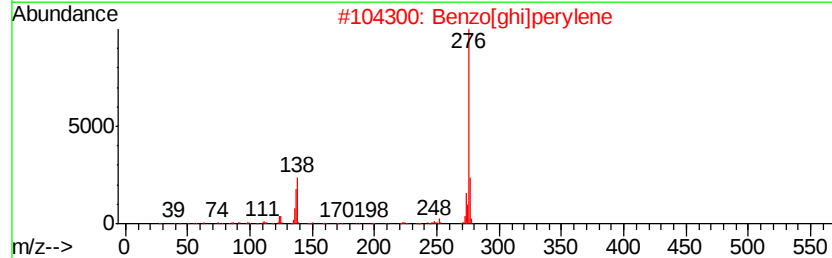
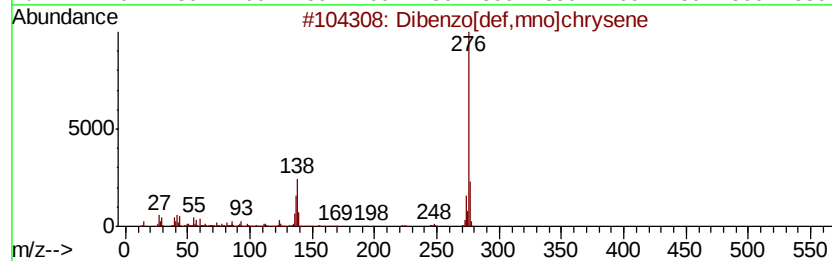
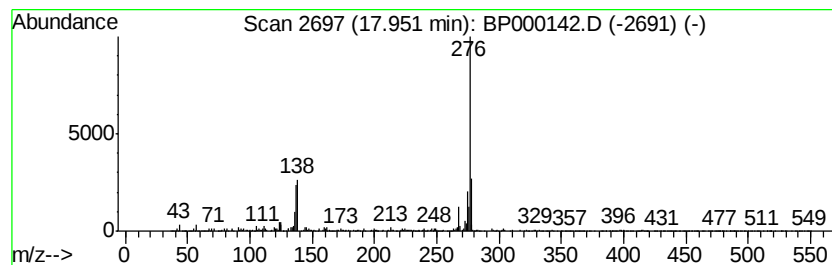
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Dibenzo[def,mno]chrysene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	3.08 ng/ul	518026	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	96
2		Benzo[ghi]perylene	276	C22H12	000191-24-2	94
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	93
4		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	91
5		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP090919\
Data File : BP000142.D
Acq On : 09 Sep 2019 19:13
Operator : HP/JU
Sample : K4633-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F2D08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.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
Butane, 2-methoxy...	2.17	176.1	ng/ul	19326100	1	6.89	2194540	20.0
n-Hexadecanoic acid	12.00	5.3	ng/ul	1092910	4	11.45	4118520	20.0
4H-Cyclopenta[def...	12.08	2.0	ng/ul	414824	4	11.45	4118520	20.0
1-Heneicosyl formate	14.02	3.9	ng/ul	1513050	5	14.15	7730030	20.0
Cyclohexane, hexa...	14.67	2.6	ng/ul	1014420	5	14.15	7730030	20.0
(DEL) Alkane: Str...	15.37	3.2	ng/ul	532370	6	15.69	3365050	20.0
(DEL) Alkane: Str...	15.80	3.4	ng/ul	572412	6	15.69	3365050	20.0
unknown-01	16.28	3.3	ng/ul	557210	6	15.69	3365050	20.0
unknown-02	16.83	3.1	ng/ul	524466	6	15.69	3365050	20.0
1,2:7,8-Dibenzoph...	17.49	11.2	ng/ul	1880500	6	15.69	3365050	20.0
Dibenzo[def,mno]c...	17.95	3.1	ng/ul	518026	6	15.69	3365050	20.0