

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081718\
 Data File : BN002370.D
 Acq On : 17 Aug 2018 09:33
 Operator : JU/SJ
 Sample : J4424-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YQ0

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.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	3.270	45	48	59	rVB	124251	189846	1.68%	0.110%
2	4.258	211	216	234	rVB	1960017	2899826	25.61%	1.682%
3	4.840	310	315	330	rVB	878168	1308225	11.55%	0.759%
4	6.828	647	653	656	rBV	988402	1759879	15.54%	1.021%
5	6.881	659	662	677	rVV	1105563	1773361	15.66%	1.029%
6	7.016	679	685	699	rVB	522726	832102	7.35%	0.483%
7	7.211	712	718	721	rBV	655854	1099843	9.71%	0.638%
8	7.246	721	724	731	rVB	611205	941959	8.32%	0.547%
9	7.675	790	797	806	rBV	616893	1017607	8.99%	0.590%
10	8.381	911	917	927	rBV	858472	1397889	12.35%	0.811%
11	8.475	927	933	947	rVB	767330	1216315	10.74%	0.706%
12	8.822	986	992	996	rBV	699969	1185581	10.47%	0.688%
13	8.863	996	999	1013	rVB	569075	930185	8.22%	0.540%
14	9.540	1108	1114	1117	rBV	571056	1036575	9.16%	0.601%
15	9.569	1117	1119	1135	rVB	530525	768457	6.79%	0.446%
16	10.075	1198	1205	1207	rBV	921550	1551863	13.71%	0.900%
17	10.446	1259	1268	1273	rBV	1068789	1908439	16.86%	1.107%
18	10.499	1273	1277	1284	rVV	862711	1389342	12.27%	0.806%
19	10.587	1285	1292	1307	rVB	613108	1150995	10.17%	0.668%
20	10.781	1318	1325	1333	rVB	873674	1403444	12.40%	0.814%
21	11.740	1481	1488	1506	rBV	1645038	2734883	24.16%	1.587%
22	12.334	1580	1589	1600	rVB	1519459	2350127	20.76%	1.363%
23	12.487	1609	1615	1622	rBV	950955	1487402	13.14%	0.863%
24	12.728	1650	1656	1663	rBV	1547378	2502750	22.11%	1.452%
25	12.787	1663	1666	1682	rVB	101455	181365	1.60%	0.105%
26	13.134	1718	1725	1729	rBV	2158437	3288365	29.04%	1.908%
27	13.175	1729	1732	1745	rVB	1040191	1505050	13.29%	0.873%
28	13.722	1818	1825	1829	rBV	1770990	2582923	22.81%	1.499%
29	13.769	1829	1833	1840	rVB	1479513	2046601	18.08%	1.187%
30	13.887	1848	1853	1861	rVB	1063862	1554072	13.73%	0.902%
31	13.998	1861	1872	1880	rBV	2108362	3066463	27.08%	1.779%
32	14.310	1918	1925	1930	rBV2	1494926	2368788	20.92%	1.374%
33	14.369	1930	1935	1945	rVB	1905579	2750245	24.29%	1.596%
34	14.528	1954	1962	1983	rBV3	1194066	2565765	22.66%	1.489%

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35	14.710	1986	1993	2003	rVB	1378946	2044768	18.06%	1.186%
36	14.940	2021	2032	2042	rBV	1161118	1681281	14.85%	0.975%
37	15.140	2056	2066	2070	rBV	578877	783441	6.92%	0.455%
38	15.175	2070	2072	2077	rVB	122026	125032	1.10%	0.073%
39	15.304	2087	2094	2099	rBV	2688590	3816296	33.71%	2.214%
40	15.357	2099	2103	2111	rVB	1264740	1705910	15.07%	0.990%
41	15.445	2111	2118	2130	rBV3	1869874	3321341	29.34%	1.927%
42	16.363	2268	2274	2282	rVB	1338491	1917879	16.94%	1.113%
43	16.710	2327	2333	2350	rBV	1786949	2636010	23.28%	1.529%
44	17.057	2385	2392	2395	rBV	2057625	2991329	26.42%	1.736%
45	17.098	2395	2399	2404	rVV	2115069	2903026	25.64%	1.684%
46	17.157	2404	2409	2412	rVV	3040056	4693812	41.46%	2.723%
47	17.186	2412	2414	2424	rVB	2390771	3075763	27.17%	1.784%
48	17.463	2453	2461	2468	rBV	1942948	2811301	24.83%	1.631%
49	17.534	2468	2473	2490	rVB	2136612	2792149	24.66%	1.620%
50	17.963	2541	2546	2550	rBV	219320	313594	2.77%	0.182%
51	18.028	2550	2557	2563	rVB	2110987	2687617	23.74%	1.559%
52	19.122	2733	2743	2752	rVB	5657132	7516981	66.39%	4.361%
53	19.245	2760	2764	2776	rBV2	86178	153160	1.35%	0.089%
54	19.457	2793	2800	2802	rBV	3929506	5698995	50.34%	3.306%
55	19.486	2802	2805	2815	rVB	4061323	5175694	45.71%	3.003%
56	20.392	2954	2959	2966	rBV	2322645	2610000	23.05%	1.514%
57	21.175	3087	3092	3097	rBV	3805942	4094887	36.17%	2.376%
58	21.239	3097	3103	3109	rVV2	6302960	11321794	100.00%	6.569%
59	21.292	3109	3112	3123	rVB	4423836	5084938	44.91%	2.950%
60	21.845	3203	3206	3210	rBV	101283	121718	1.08%	0.071%
61	22.304	3281	3284	3289	rVB	175509	207234	1.83%	0.120%
62	22.810	3363	3370	3383	rVB	3737093	6035116	53.31%	3.501%
63	23.339	3451	3460	3463	rBV2	3223715	6331866	55.93%	3.674%
64	23.380	3463	3467	3475	rVV	2955487	5408571	47.77%	3.138%
65	23.474	3475	3483	3494	rVB	2097371	3976470	35.12%	2.307%
66	23.998	3566	3572	3580	rVB	249372	494376	4.37%	0.287%
67	24.727	3690	3696	3704	rBV	155165	346010	3.06%	0.201%
68	25.586	3837	3842	3848	rVB3	81729	183937	1.62%	0.107%
69	25.704	3849	3862	3877	rVV2	3760176	10551172	93.19%	6.122%

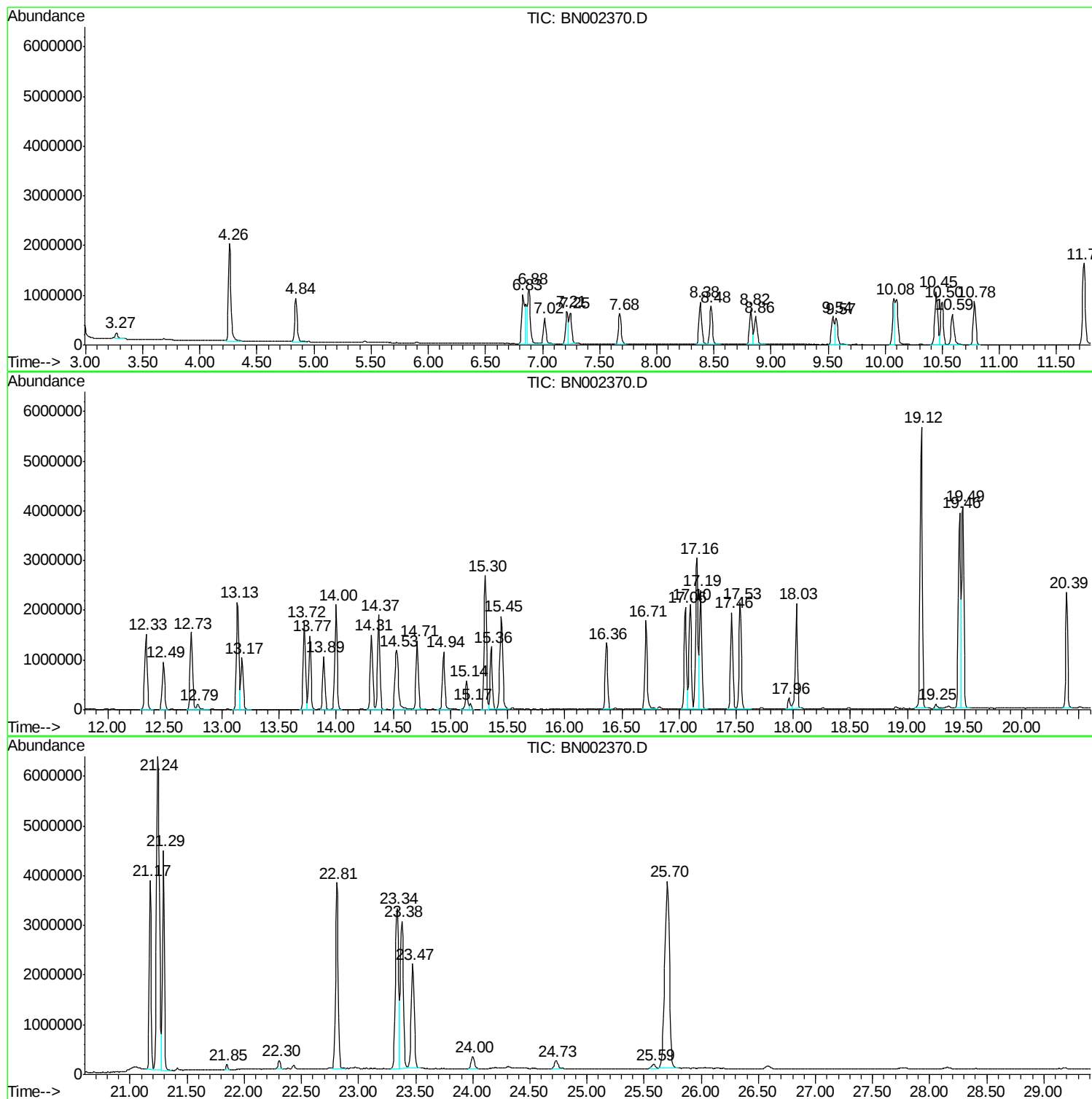
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Quant Title : SVOA CALIBRATION

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



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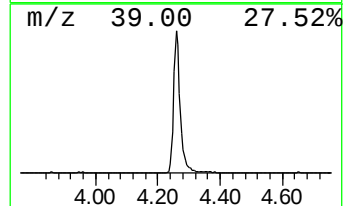
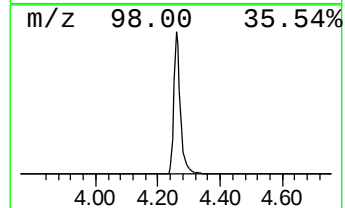
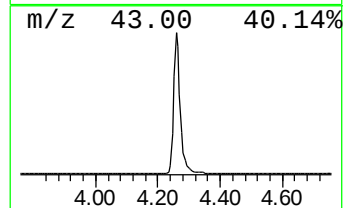
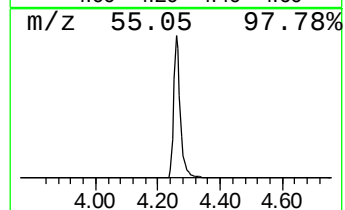
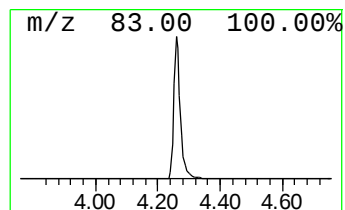
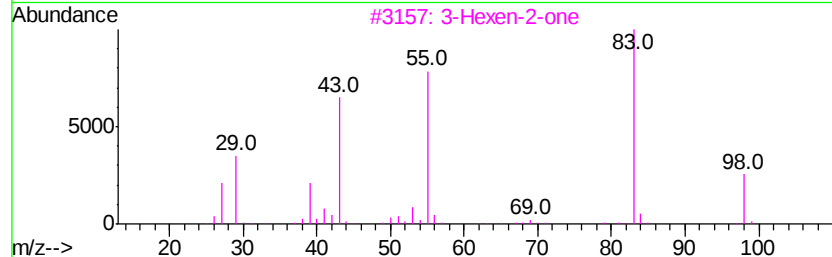
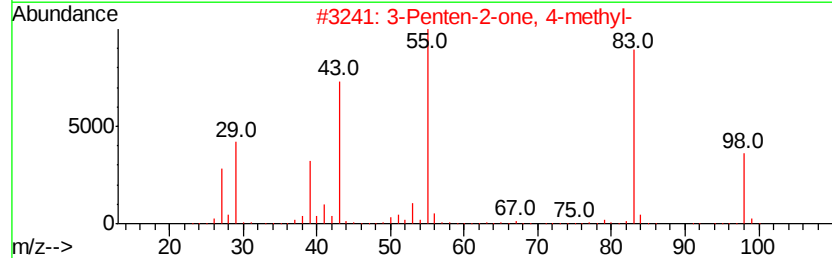
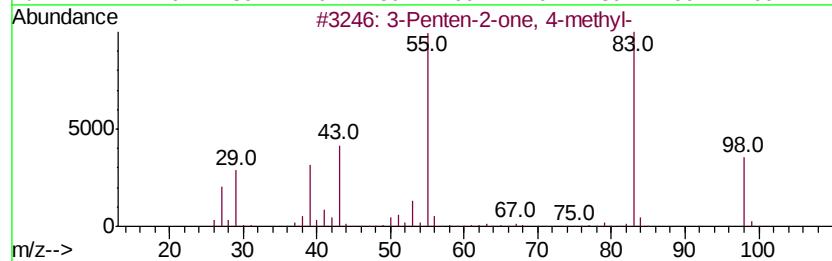
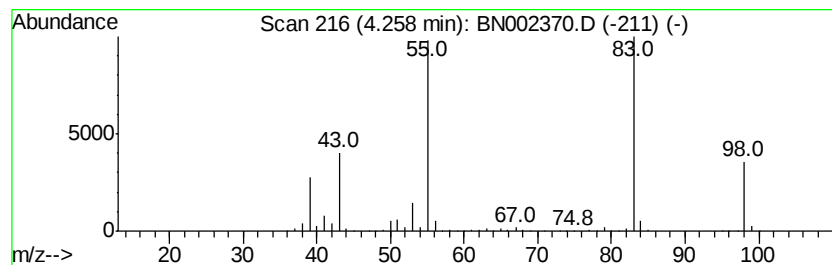
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.26	56.99 ng/ul	2899830	1,4-Dichlorobenzene-d4	7.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3			3-Hexen-2-one	98	C6H10O	000763-93-9	91
4			3-Hexen-2-one	98	C6H10O	000763-93-9	90
5			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86



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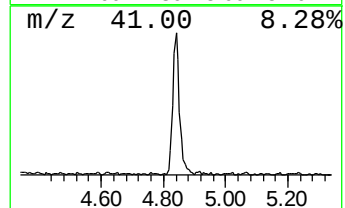
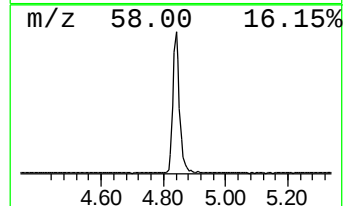
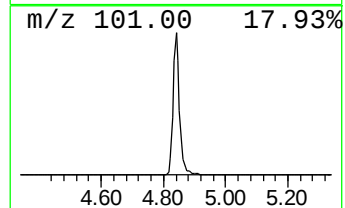
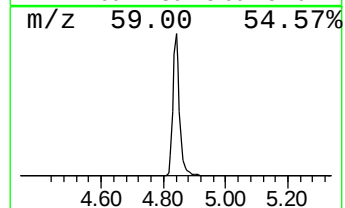
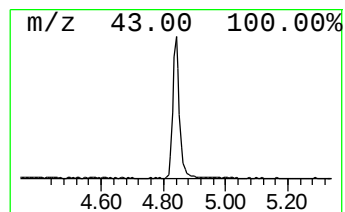
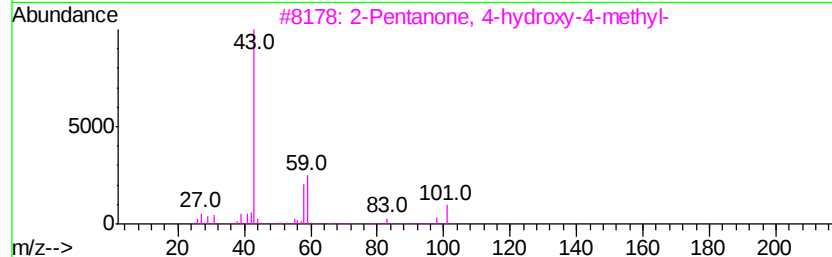
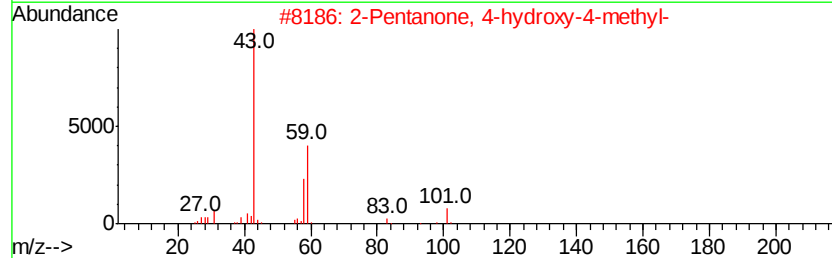
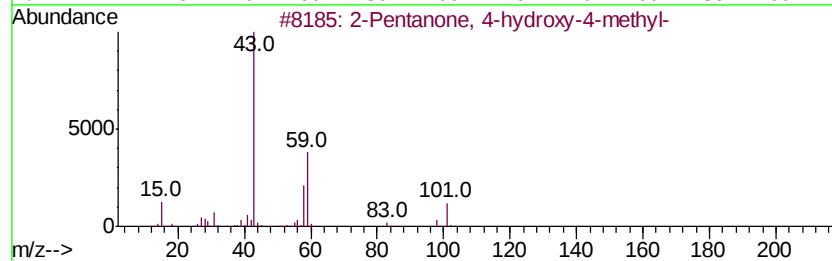
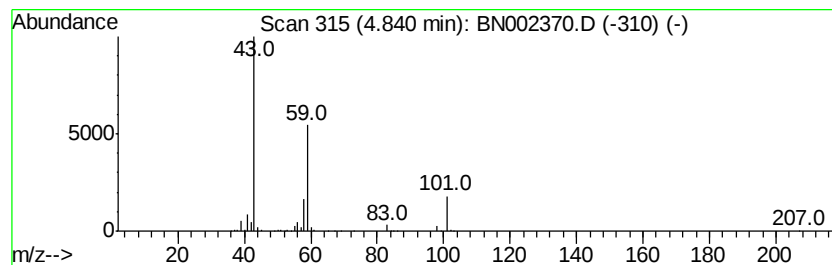
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	25.71 ng/ul	1308230	1,4-Dichlorobenzene-d4	7.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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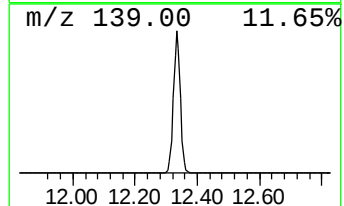
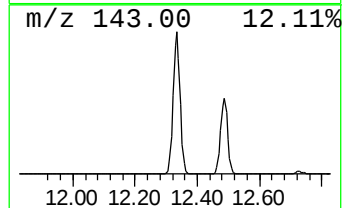
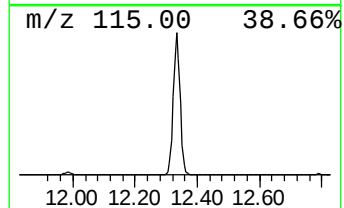
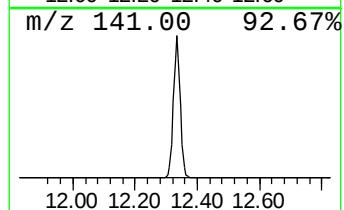
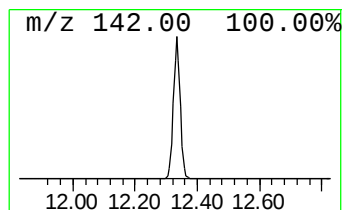
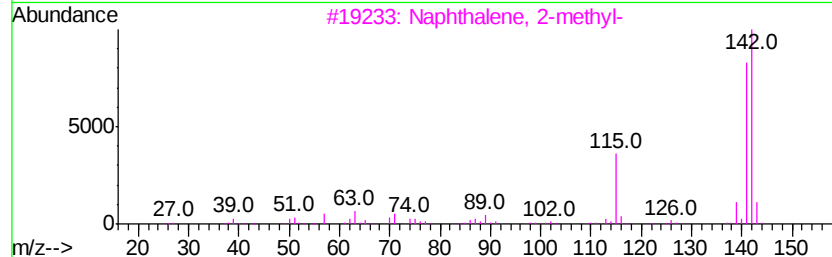
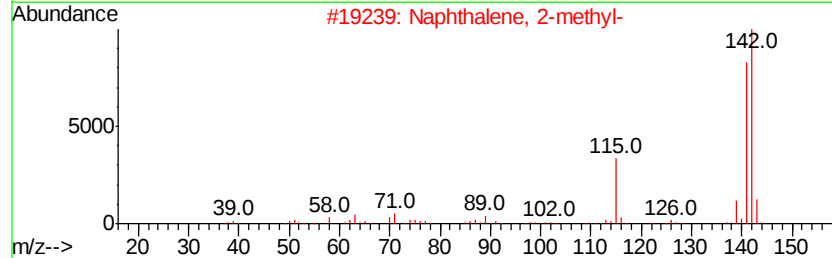
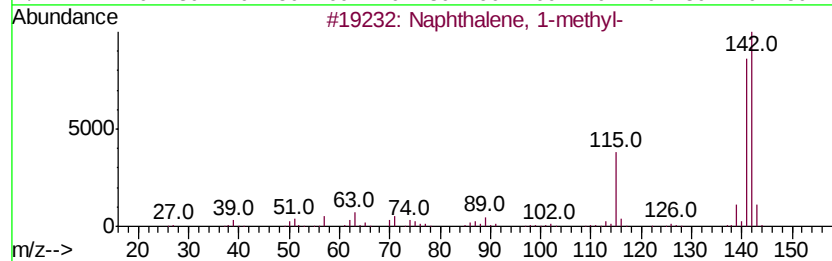
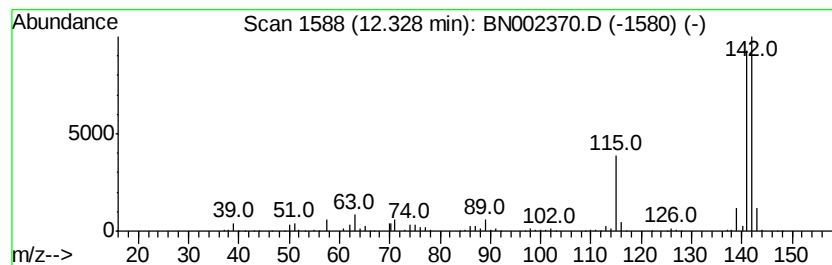
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Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	24.63 ng/ul	2350130	Naphthalene-d8	10.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91



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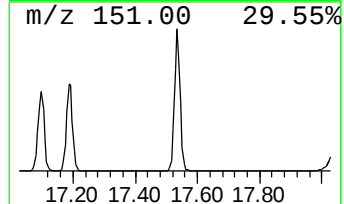
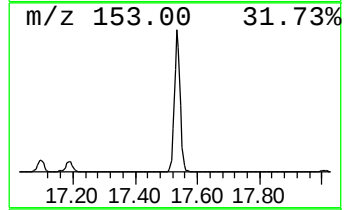
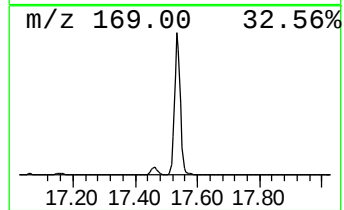
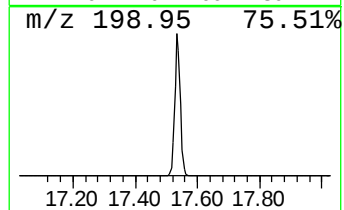
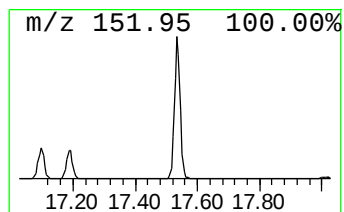
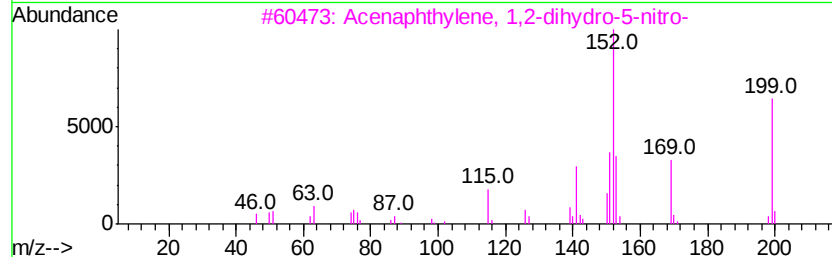
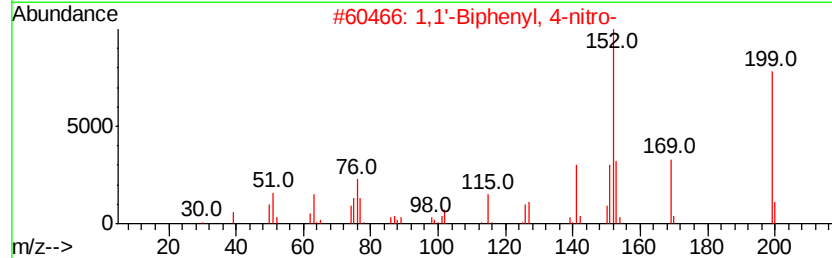
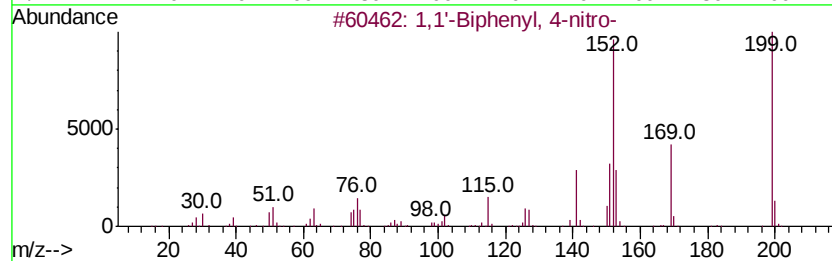
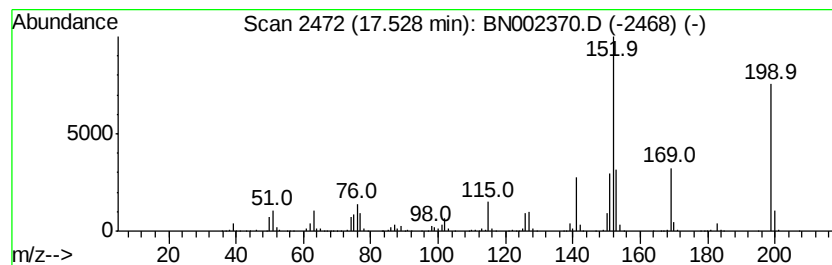
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Peak Number 4 1,1'-Biphenyl, 4-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.53	18.67 ng/ul	2792150	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	99
2		1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3		Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	93
4		1,1'-Biphenyl, 3-nitro-	199	C12H9NO2	002113-58-8	90
5		Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081718\
Data File : BN002370.D
Acq On : 17 Aug 2018 09:33
Operator : JU/SJ
Sample : J4424-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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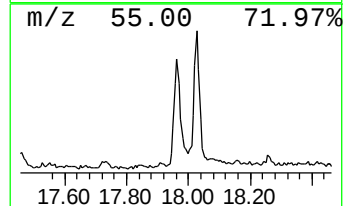
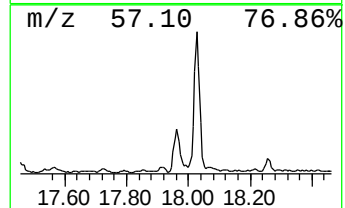
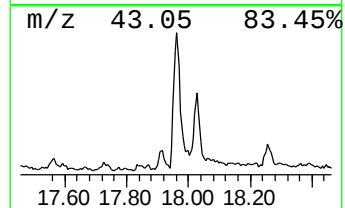
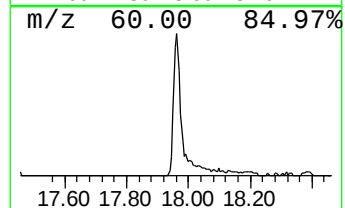
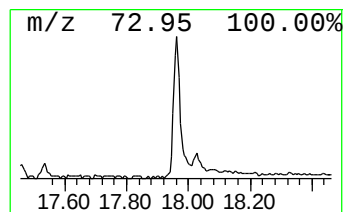
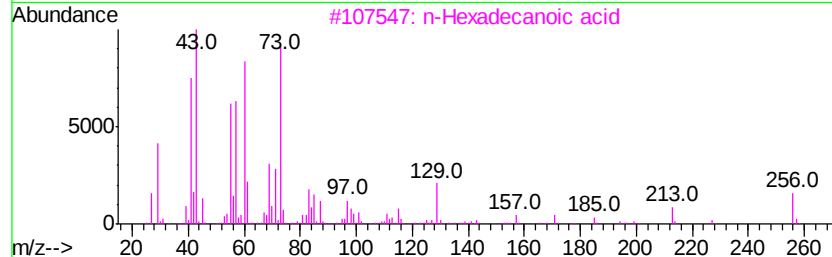
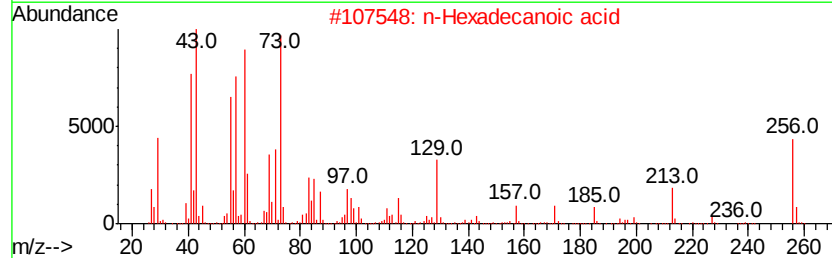
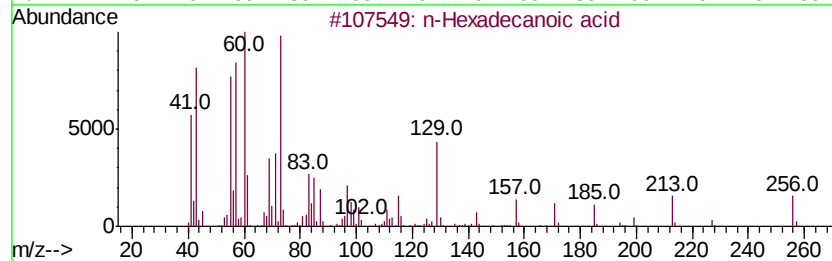
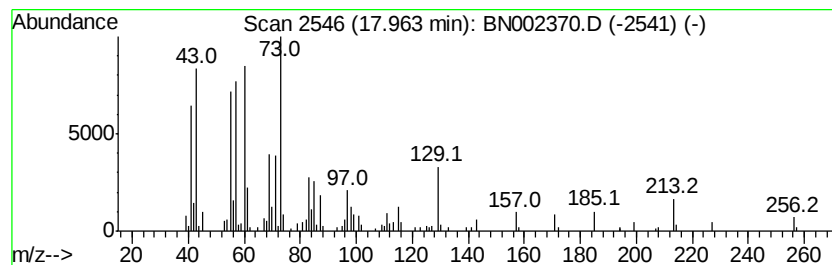
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	2.10 ng/ul	313594	Phenanthrene-d10	17.06

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	98	
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96	
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	90	
5	Tridecanoic acid	214	C13H26O2	000638-53-9	87	



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081718\
Data File : BN002370.D
Acq On : 17 Aug 2018 09:33
Operator : JU/SJ
Sample : J4424-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3YQ0

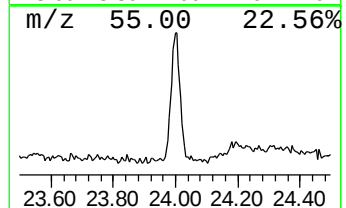
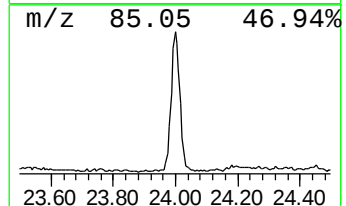
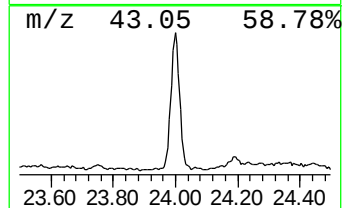
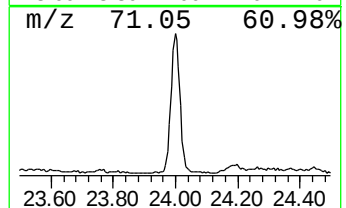
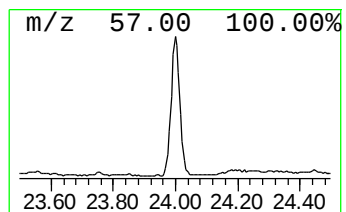
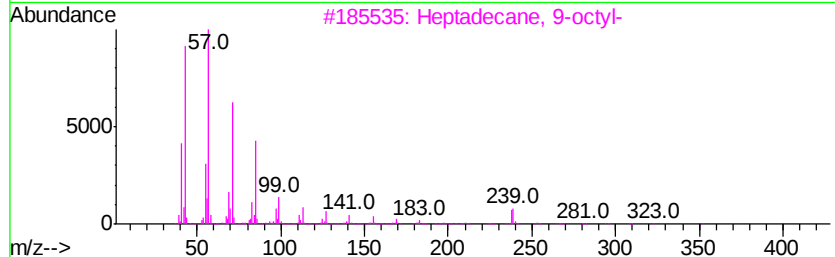
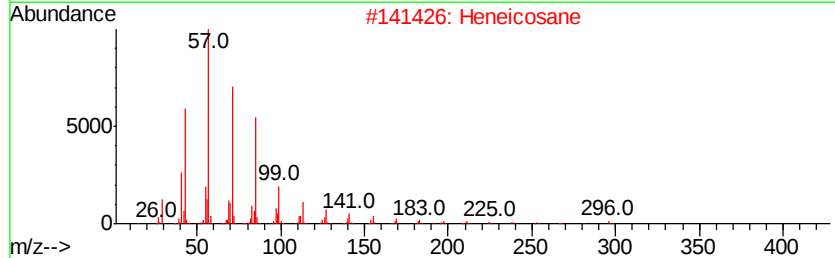
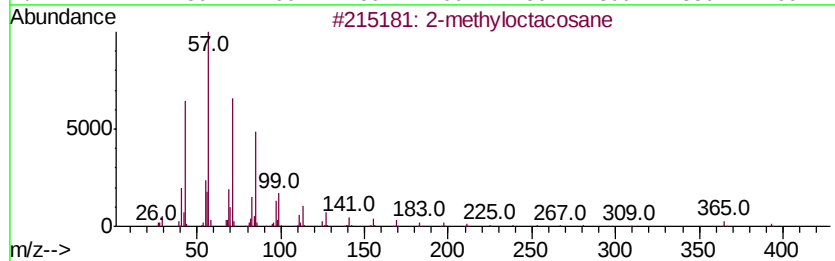
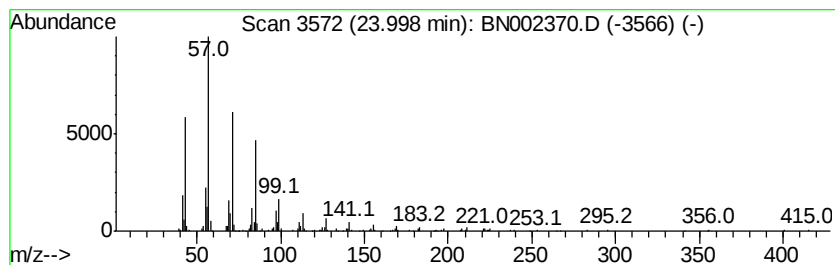
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	2.49 ng/ul	494376	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		Tetracosane	338	C24H50	000646-31-1	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081718\
Data File : BN002370.D
Acq On : 17 Aug 2018 09:33
Operator : JU/SJ
Sample : J4424-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3YQ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.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
3-Penten-2-one, 4...	4.26	57.0	ng/ul	2899830	1	7.68	1017610	20.0
2-Pentanone, 4-hy...	4.84	25.7	ng/ul	1308230	1	7.68	1017610	20.0
Naphthalene, 1-me...	12.33	24.6	ng/ul	2350130	2	10.45	1908440	20.0
1,1'-Biphenyl, 4-...	17.53	18.7	ng/ul	2792150	4	17.06	2991330	20.0
n-Hexadecanoic acid	17.96	2.1	ng/ul	313594	4	17.06	2991330	20.0
(DEL) Alkane: Str...	24.00	2.5	ng/ul	494376	6	23.47	3976470	20.0