

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
 Data File : BP000173.D
 Acq On : 10 Sep 2019 18:36
 Operator : HP/JU
 Sample : K4780-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0BM4

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	2.163	7	13	23	rBV	3839850	4817976	69.18%	5.665%
2	2.610	85	89	98	rBV	74187	109145	1.57%	0.128%
3	3.922	306	312	316	rBV2	8057	11793	0.17%	0.014%
4	3.969	316	320	323	rVV	23193	33290	0.48%	0.039%
5	3.999	323	325	330	rVB	21485	25300	0.36%	0.030%
6	4.193	354	358	363	rBV	16335	20124	0.29%	0.024%
7	4.799	457	461	466	rVB	8686	9762	0.14%	0.011%
8	5.087	507	510	517	rBV2	8157	12835	0.18%	0.015%
9	6.022	665	669	672	rVB	25444	20131	0.29%	0.024%
10	6.198	693	699	705	rVB2	18831	32607	0.47%	0.038%
11	6.440	730	740	747	rBV4	8960	17227	0.25%	0.020%
12	6.504	747	751	758	rBV	795212	645256	9.27%	0.759%
13	6.569	758	762	766	rBV	2164800	1834360	26.34%	2.157%
14	6.657	773	777	780	rBV	2575894	2013957	28.92%	2.368%
15	6.893	813	817	820	rBV	3151244	2414017	34.66%	2.839%
16	6.951	824	827	831	rBV	11314	8392	0.12%	0.010%
17	7.257	875	879	882	rBV	2081275	1655200	23.77%	1.946%
18	7.445	907	911	914	rBV	2775648	2150684	30.88%	2.529%
19	7.528	921	925	931	rVB3	9386	11542	0.17%	0.014%
20	7.663	945	948	951	rVB	38713	30699	0.44%	0.036%
21	7.775	963	967	970	rBV	2114902	1703775	24.46%	2.003%
22	7.951	994	997	1000	rBV	36533	27620	0.40%	0.032%
23	8.010	1004	1007	1011	rBV	3124950	2767112	39.73%	3.254%
24	8.175	1032	1035	1039	rVB	4124296	3297301	47.34%	3.877%
25	8.228	1041	1044	1047	rBV	759222	552490	7.93%	0.650%
26	8.498	1085	1090	1095	rBV5	8556	11935	0.17%	0.014%
27	8.716	1123	1127	1131	rBV	105852	86397	1.24%	0.102%
28	8.857	1148	1151	1156	rBV	66985	53907	0.77%	0.063%
29	9.134	1195	1198	1201	rBV	11354	13410	0.19%	0.016%
30	9.251	1214	1218	1221	rBV2	27935	34960	0.50%	0.041%
31	9.334	1228	1232	1233	rBV	9455	9315	0.13%	0.011%
32	9.351	1233	1235	1238	rVB2	10542	11455	0.16%	0.013%
33	9.416	1243	1246	1253	rVB2	11320	16158	0.23%	0.019%
34	9.634	1279	1283	1285	rBV	4961975	4276804	61.41%	5.029%

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 Title : SVOA CALIBRATION

35	9.651	1285	1286	1292	rVB	644358	375777	5.40%	0.442%
36	9.751	1301	1303	1305	rBV3	7048	7916	0.11%	0.009%
37	9.787	1305	1309	1313	rBV	6162977	5289436	75.95%	6.220%
38	9.881	1322	1325	1329	rVB2	13797	11840	0.17%	0.014%
39	9.945	1332	1336	1339	rBV	5238962	4126439	59.25%	4.852%
40	10.022	1346	1349	1357	rBV	483416	480864	6.90%	0.565%
41	10.145	1364	1370	1376	rVB4	28802	44932	0.65%	0.053%
42	10.245	1383	1387	1390	rBV4	9634	11342	0.16%	0.013%
43	10.322	1398	1400	1403	rVB	19569	15351	0.22%	0.018%
44	10.363	1403	1407	1409	rVV	28564	25176	0.36%	0.030%
45	10.387	1409	1411	1413	rVB	16778	11039	0.16%	0.013%
46	10.469	1421	1425	1428	rBV	7776102	6238149	89.57%	7.335%
47	10.528	1431	1435	1444	rVV	1463087	1393435	20.01%	1.639%
48	10.598	1444	1447	1451	rVB	99931	87225	1.25%	0.103%
49	10.781	1474	1478	1480	rBV2	8484	9262	0.13%	0.011%
50	10.839	1485	1488	1490	rVV	19327	22205	0.32%	0.026%
51	10.869	1490	1493	1495	rVV	19253	22208	0.32%	0.026%
52	10.898	1495	1498	1501	rVB2	42248	40188	0.58%	0.047%
53	11.263	1557	1560	1563	rVB	18126	14553	0.21%	0.017%
54	11.322	1566	1570	1575	rBV	31386	32412	0.47%	0.038%
55	11.410	1581	1585	1587	rBV4	5022	7856	0.11%	0.009%
56	11.445	1587	1591	1595	rBV	4812225	4285111	61.53%	5.039%
57	11.504	1597	1601	1605	rBV	7428471	6468215	92.88%	7.606%
58	11.622	1618	1621	1625	rVB	21755	20777	0.30%	0.024%
59	11.657	1625	1627	1633	rBV2	7572	8340	0.12%	0.010%
60	11.757	1641	1644	1650	rVB3	14286	13633	0.20%	0.016%
61	11.992	1680	1684	1703	rBV2	560479	663107	9.52%	0.780%
62	12.175	1710	1715	1718	rVB2	28417	34527	0.50%	0.041%
63	12.575	1780	1783	1787	rVB	46559	40666	0.58%	0.048%
64	12.616	1787	1790	1794	rBV7	10582	12442	0.18%	0.015%
65	12.663	1794	1798	1801	rBV	81376	80414	1.15%	0.095%
66	12.698	1801	1804	1806	rBV	34308	39503	0.57%	0.046%
67	12.798	1817	1821	1823	rBV	230894	234747	3.37%	0.276%
68	12.822	1823	1825	1833	rVV	104502	124457	1.79%	0.146%
69	12.898	1833	1838	1842	rVV	7563062	6964420	100.00%	8.189%
70	12.957	1845	1848	1850	rVV	92962	94508	1.36%	0.111%
71	12.980	1850	1852	1860	rVV2	58997	59081	0.85%	0.069%

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72	13.063	1863	1866	1869	rBV2	12576	12589	0.18%	0.015%
73	13.328	1905	1911	1918	rBV	144725	183716	2.64%	0.216%
74	13.551	1944	1949	1951	rBV5	30311	43365	0.62%	0.051%
75	13.651	1963	1966	1967	rBV2	11723	11313	0.16%	0.013%
76	13.680	1967	1971	1980	rVB	240910	290270	4.17%	0.341%
77	13.933	2012	2014	2020	rVB5	15783	16475	0.24%	0.019%
78	14.022	2022	2029	2045	rBV	739883	1185143	17.02%	1.394%
79	14.139	2045	2049	2053	rBV	4073152	3855349	55.36%	4.533%
80	14.351	2081	2085	2095	rVB	356221	472290	6.78%	0.555%
81	14.669	2134	2139	2149	rBV	347313	533157	7.66%	0.627%
82	15.004	2191	2196	2203	rBV	336790	526454	7.56%	0.619%
83	15.174	2222	2225	2229	rVB	68819	82474	1.18%	0.097%
84	15.380	2254	2260	2267	rBV	289870	497983	7.15%	0.586%
85	15.586	2288	2295	2304	rBV	4220094	5841477	83.88%	6.869%
86	15.686	2305	2312	2323	rVB	2559279	3662023	52.58%	4.306%
87	15.798	2325	2331	2340	rVB	256590	526057	7.55%	0.619%
88	16.280	2407	2413	2421	rBV	220446	503306	7.23%	0.592%
89	16.833	2502	2507	2524	rVB	157539	410274	5.89%	0.482%
90	17.492	2613	2619	2631	rVB2	89383	241062	3.46%	0.283%

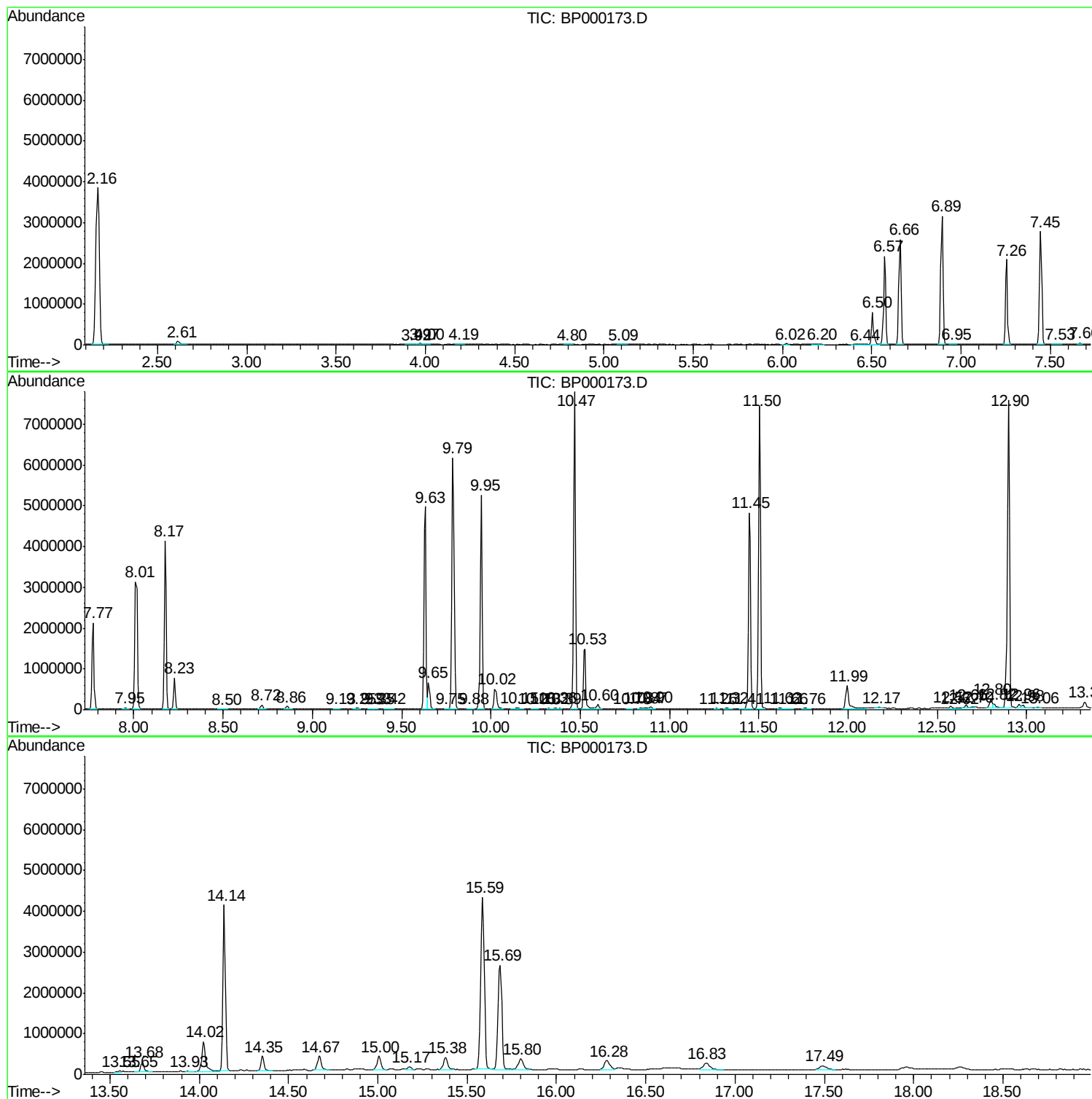
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 10 Sample Multiplier: 1

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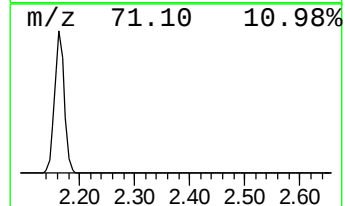
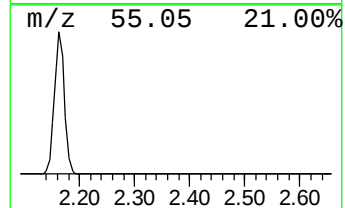
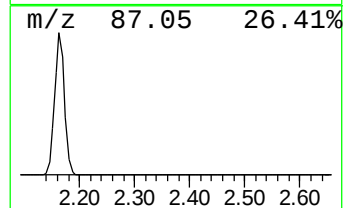
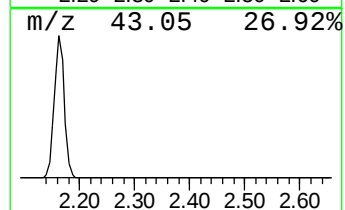
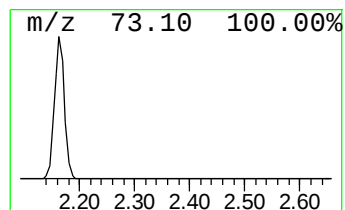
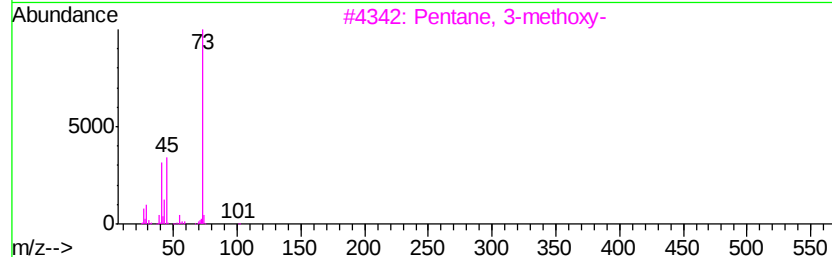
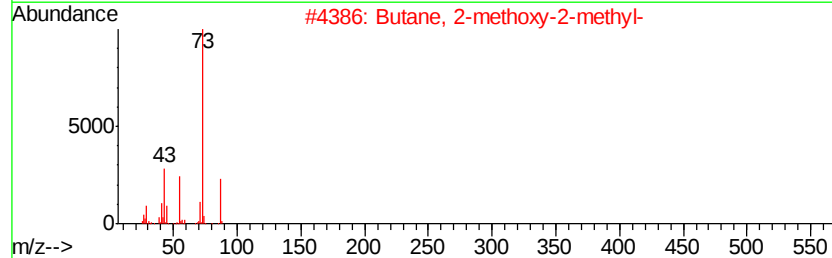
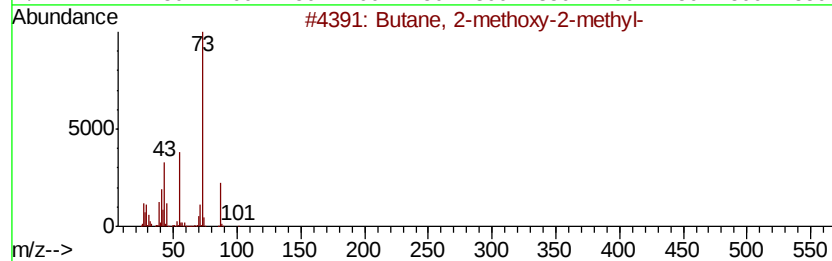
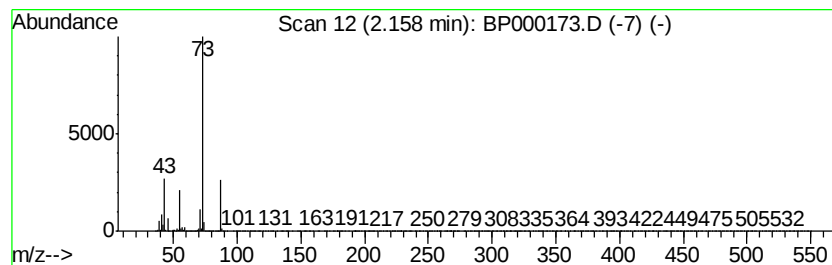
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION

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.
2.16	39.92 ng/ul	4817980	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	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	7



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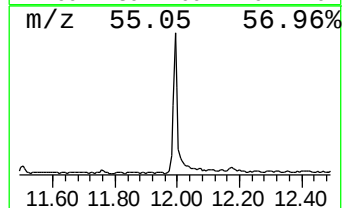
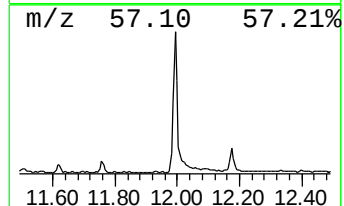
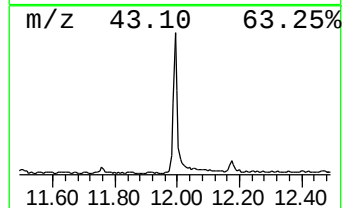
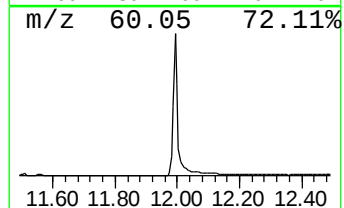
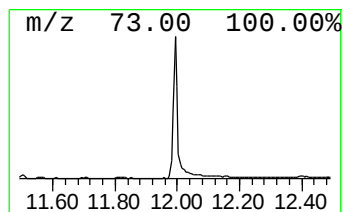
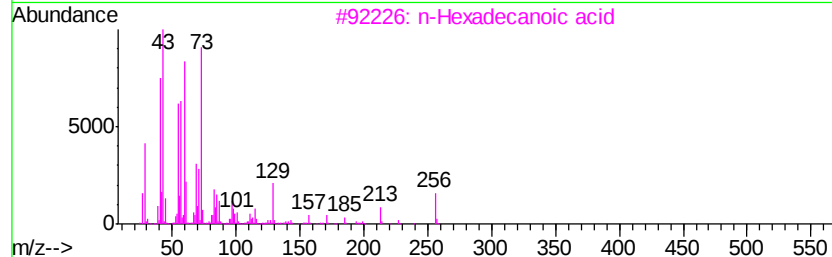
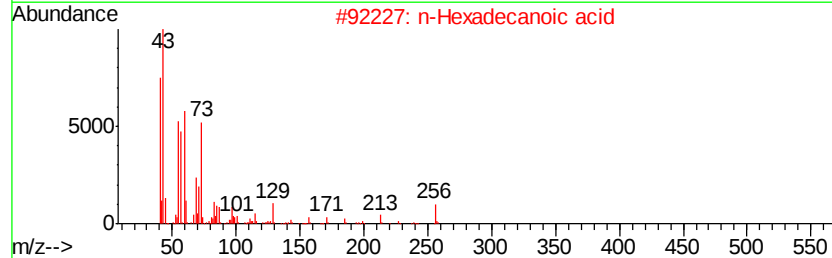
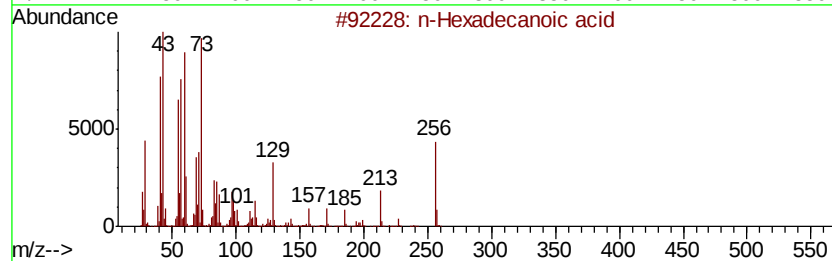
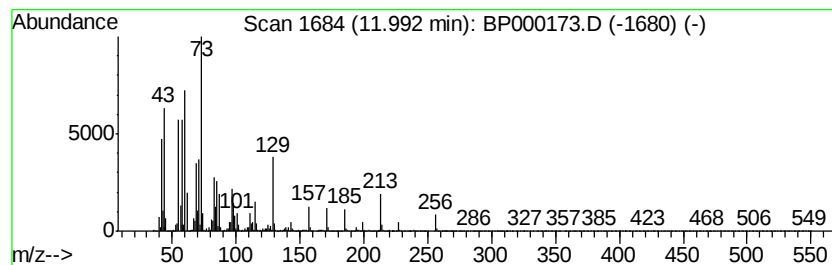
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Quant Title : SVOA CALIBRATION

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.	
11.99	3.09 ng/ul	663107	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	90
4	Tridecanoic acid		214	C13H26O2	000638-53-9	89
5	Pentadecanoic acid		242	C15H30O2	001002-84-2	74



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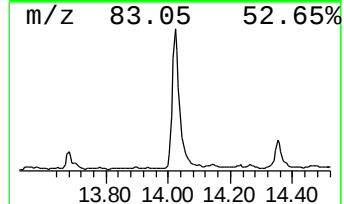
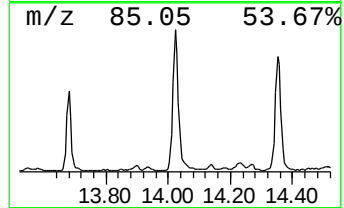
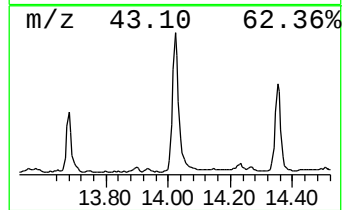
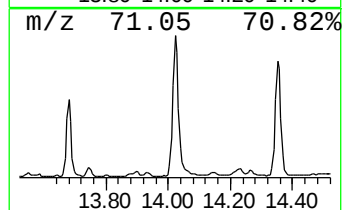
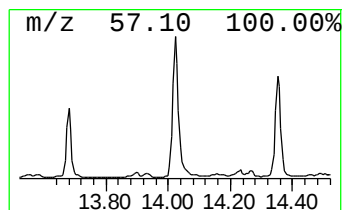
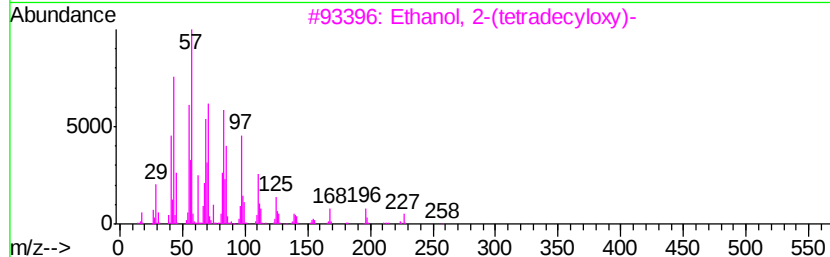
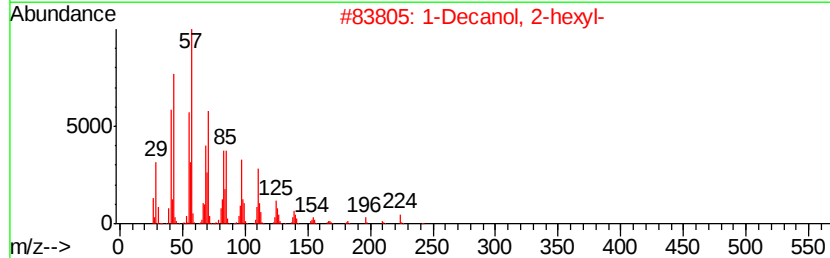
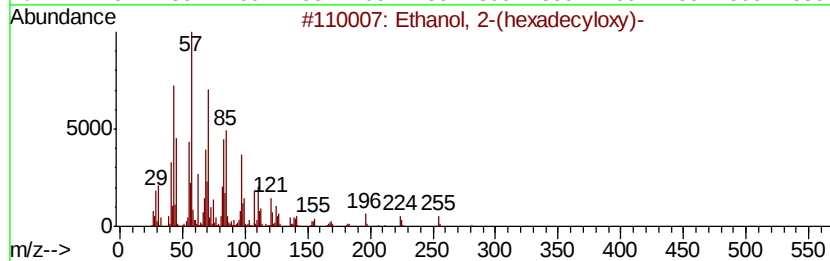
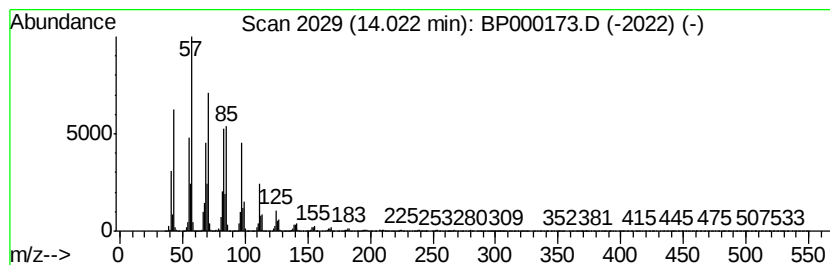
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(hexadecyloxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	6.15 ng/ul	1185140	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	91
2		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	90
3		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
4		Eicosane	282	C20H42	000112-95-8	86
5		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	86



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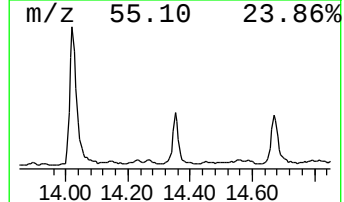
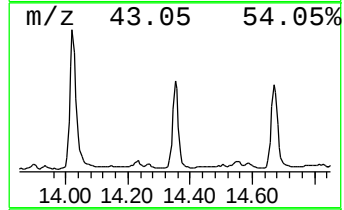
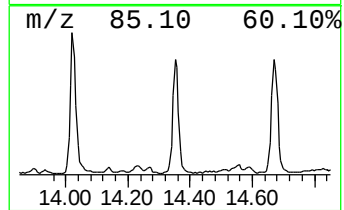
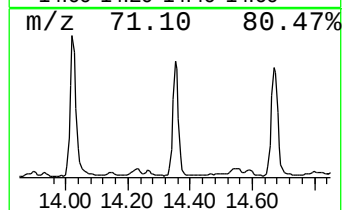
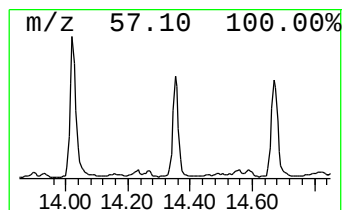
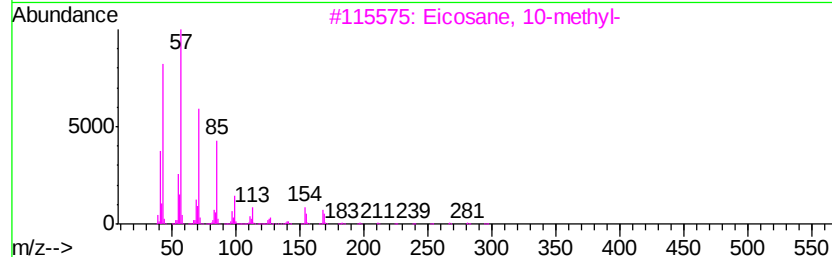
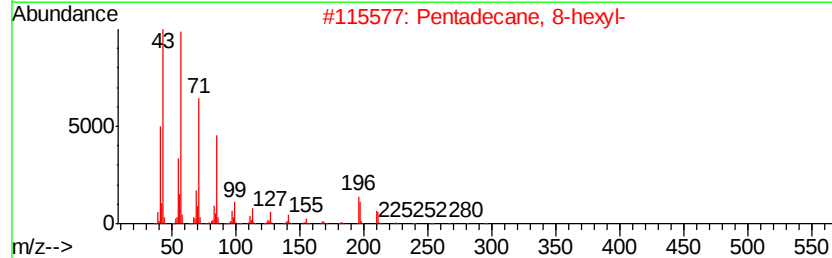
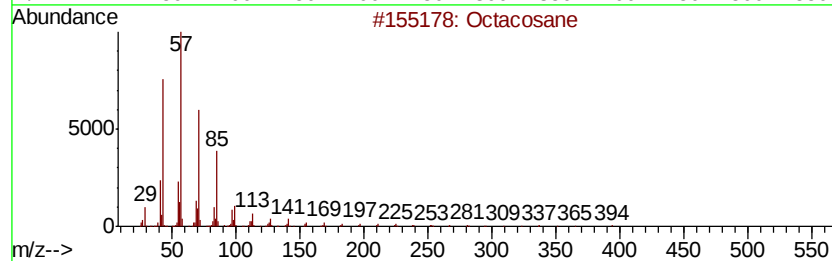
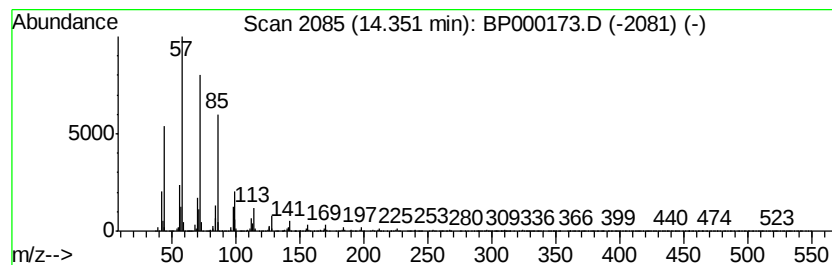
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	2.45 ng/ul	472290	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	98
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0BM4

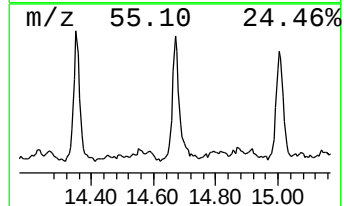
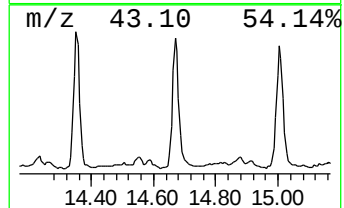
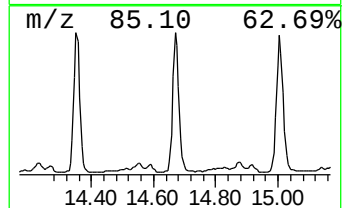
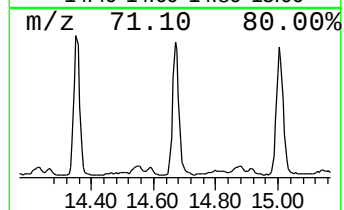
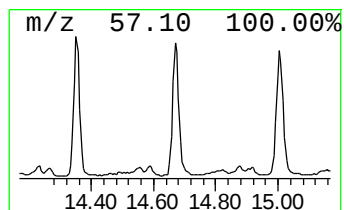
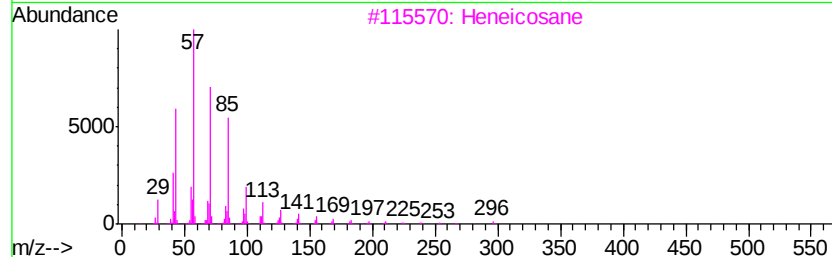
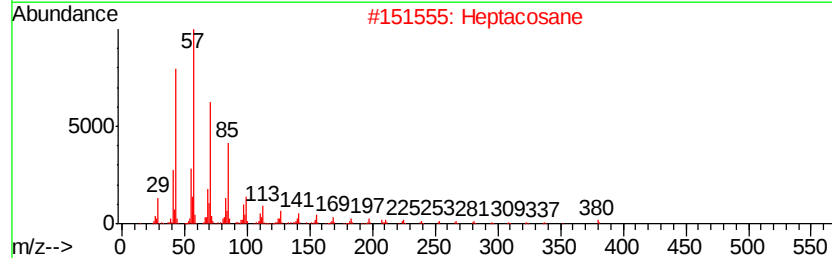
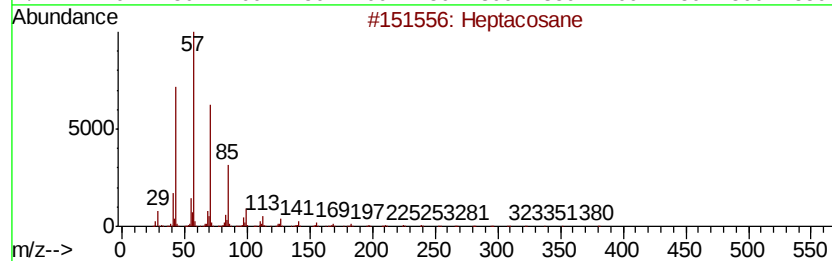
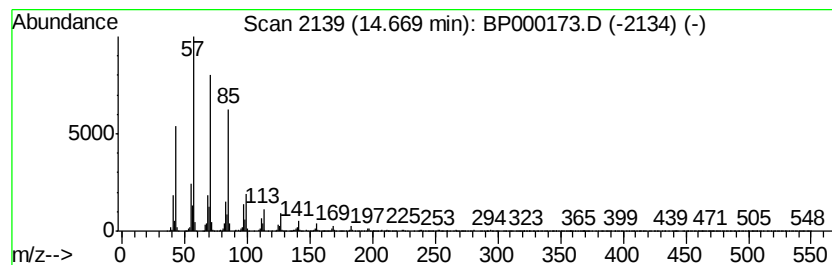
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.77 ng/ul	533157	Chrysene-d12	14.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	94
2		Heptacosane	380	C27H56	000593-49-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0BM4

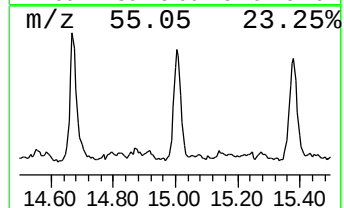
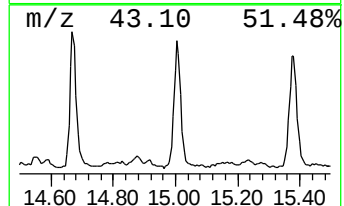
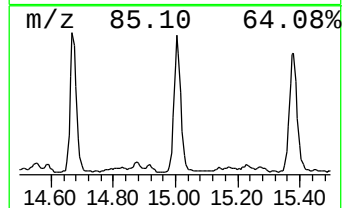
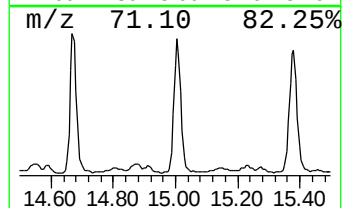
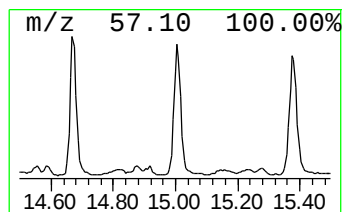
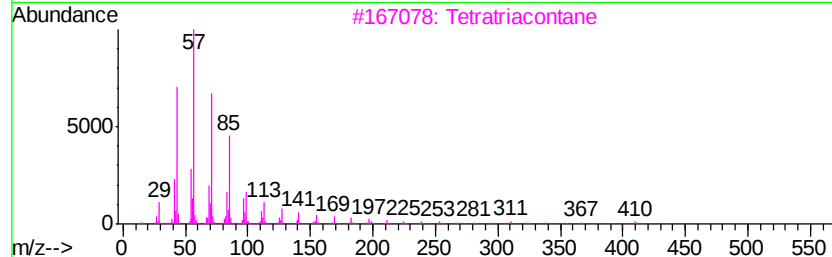
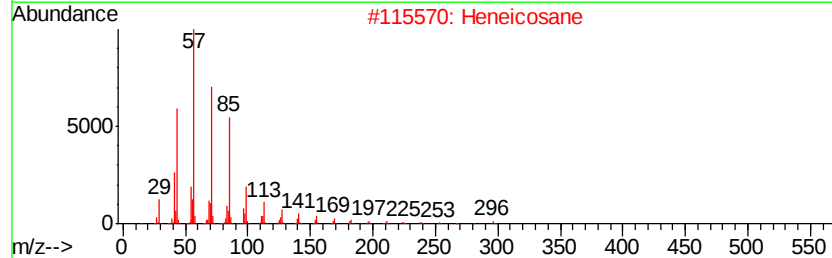
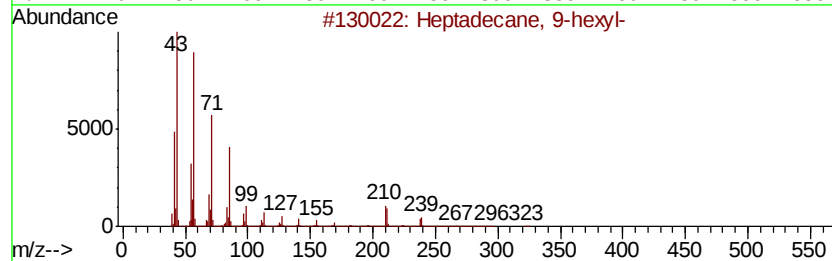
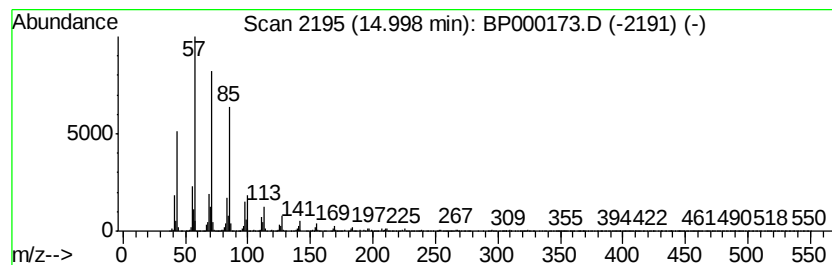
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	2.88 ng/ul	526454	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetratriacontane	479	C34H70	014167-59-0	87
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	87
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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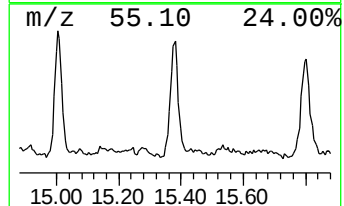
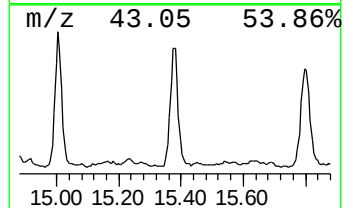
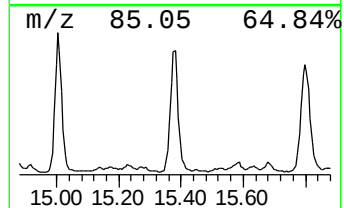
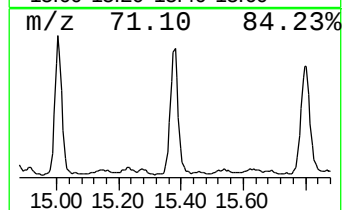
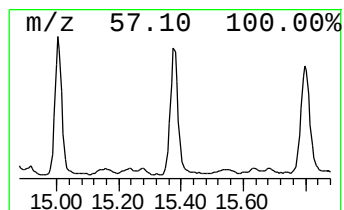
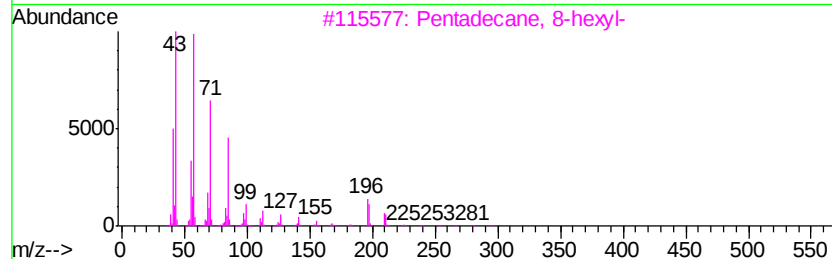
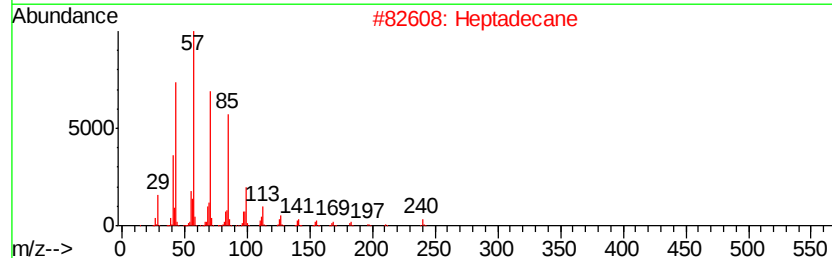
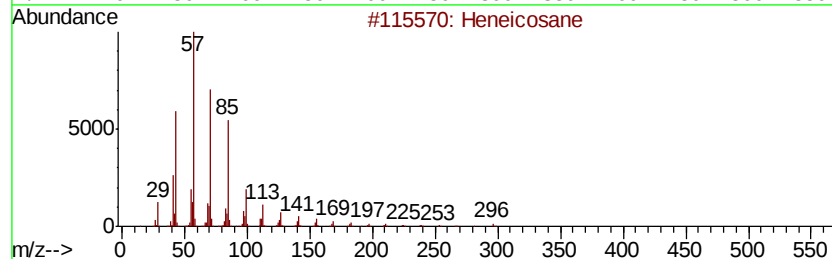
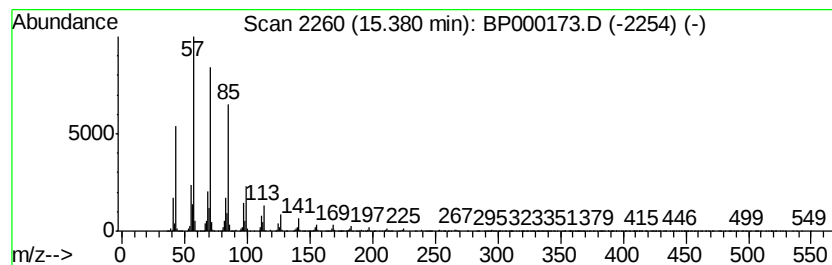
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.38	2.72 ng/ul	497983	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
4		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	90
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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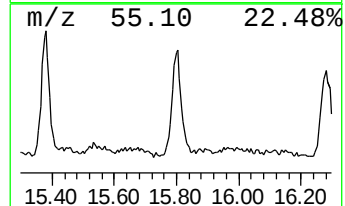
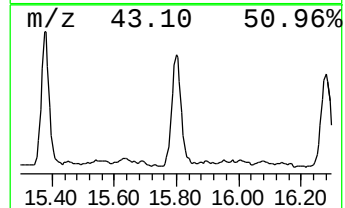
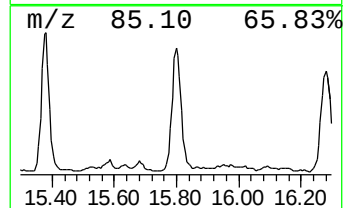
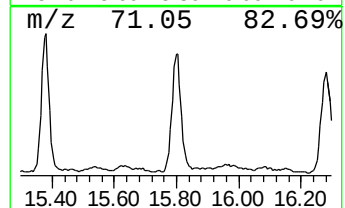
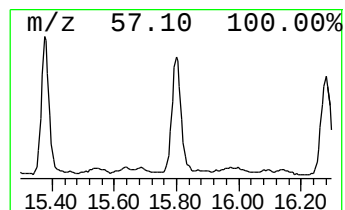
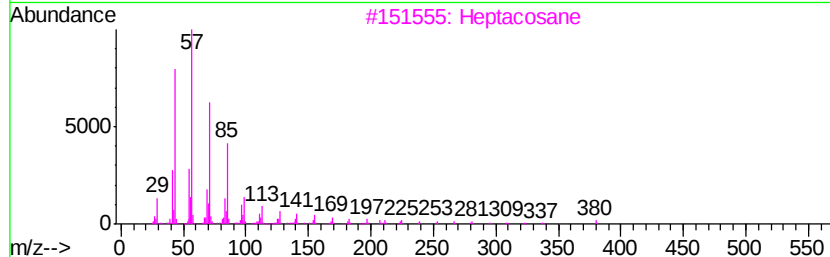
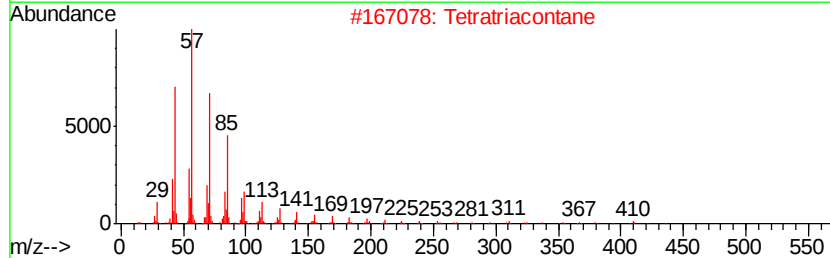
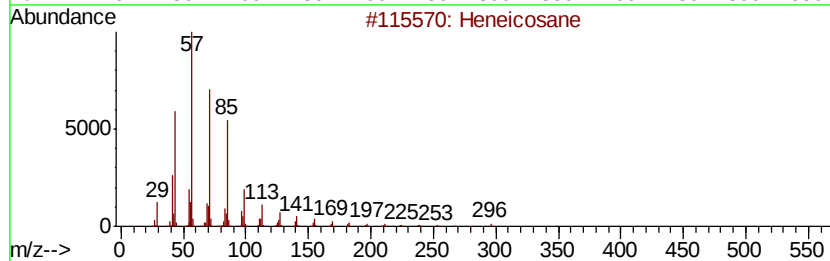
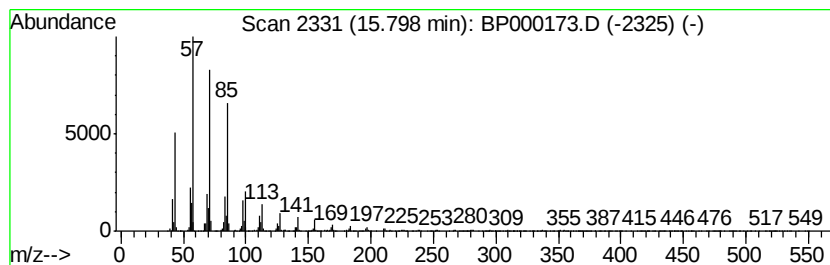
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	2.87 ng/ul	526057	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Tetratriacontane	479	C34H70	014167-59-0	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0BM4

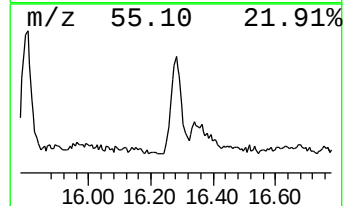
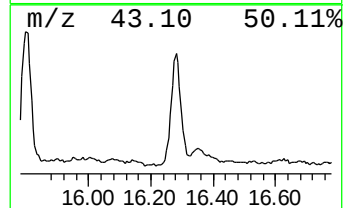
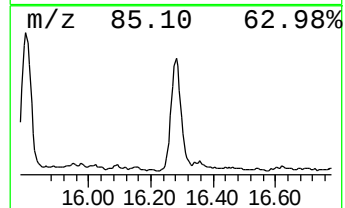
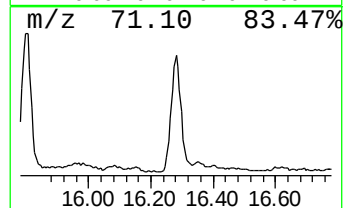
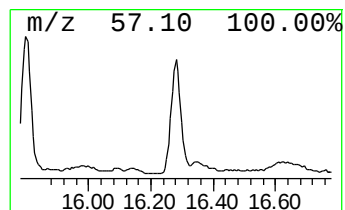
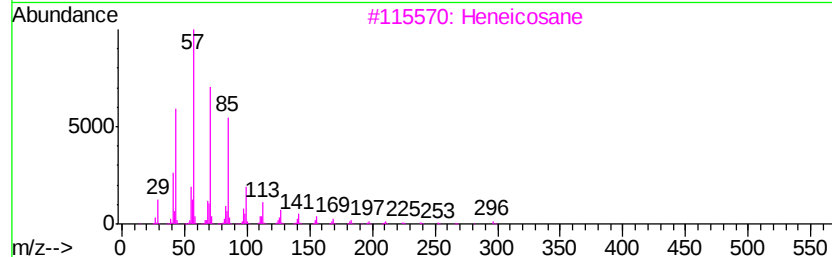
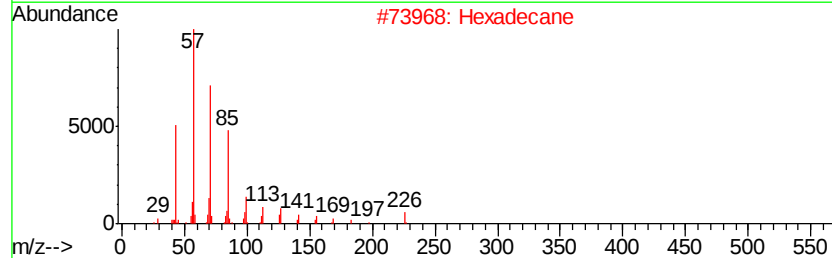
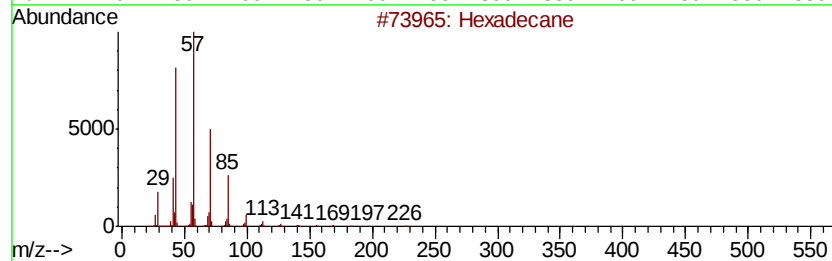
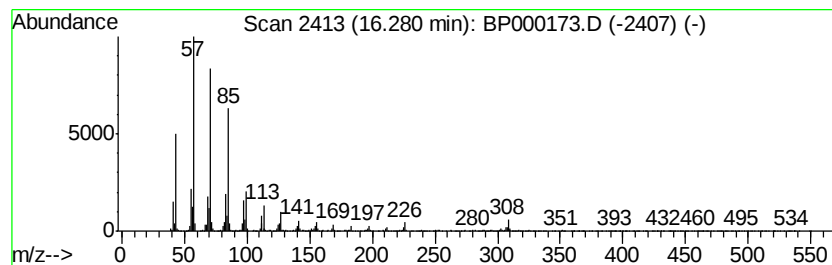
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	2.75 ng/ul	503306	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	95
2		Hexadecane	226	C16H34	000544-76-3	95
3		Heneicosane	296	C21H44	000629-94-7	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
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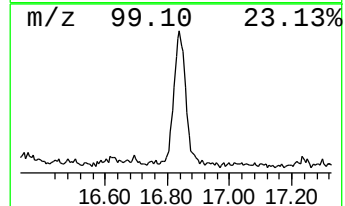
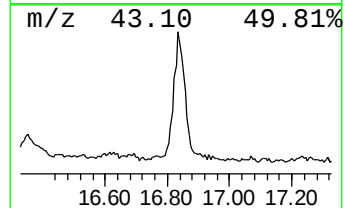
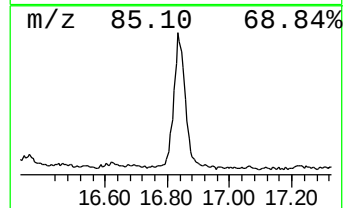
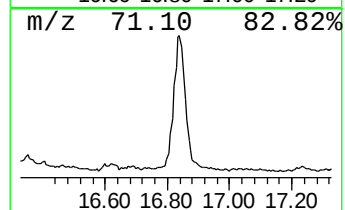
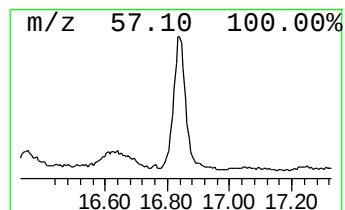
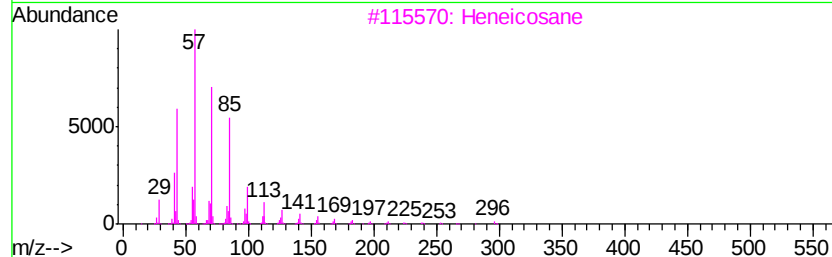
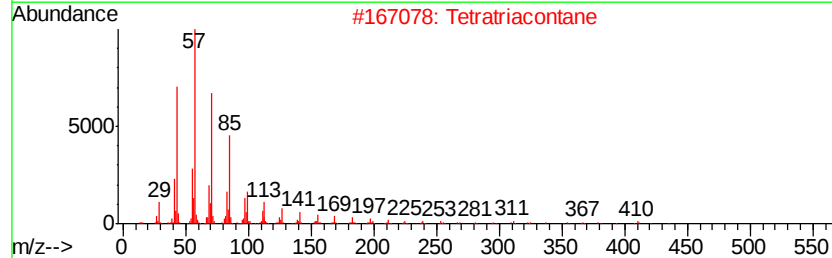
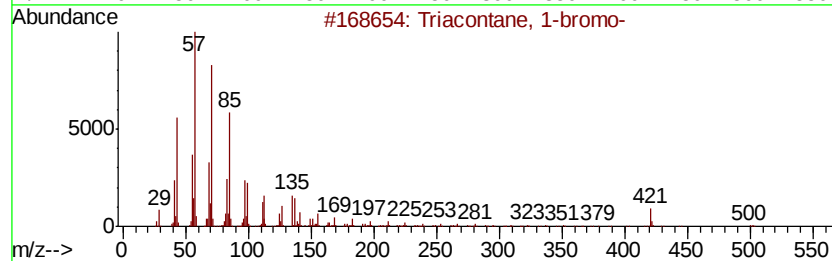
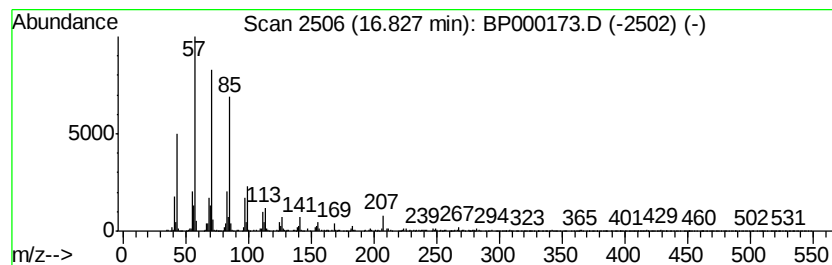
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 Quant Title : SVOA CALIBRATION

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

 Peak Number 10 Triacontane, 1-bromo- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.83	2.24 ng/ul	410274	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	95
2		Tetratriacontane	479	C34H70	014167-59-0	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
Data File : BP000173.D
Acq On : 10 Sep 2019 18:36
Operator : HP/JU
Sample : K4780-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0BM4

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...	2.16	39.9	ng/ul	4817980	1	6.89	2414020	20.0
n-Hexadecanoic acid	11.99	3.1	ng/ul	663107	4	11.45	4285110	20.0
Ethanol, 2-(hexad...	14.02	6.2	ng/ul	1185140	5	14.14	3855350	20.0
(DEL) Alkane: Str...	14.35	2.5	ng/ul	472290	5	14.14	3855350	20.0
(DEL) Alkane: Str...	14.67	2.8	ng/ul	533157	5	14.14	3855350	20.0
(DEL) Alkane: Str...	15.00	2.9	ng/ul	526454	6	15.69	3662020	20.0
(DEL) Alkane: Str...	15.38	2.7	ng/ul	497983	6	15.69	3662020	20.0
(DEL) Alkane: Str...	15.80	2.9	ng/ul	526057	6	15.69	3662020	20.0
(DEL) Alkane: Str...	16.28	2.8	ng/ul	503306	6	15.69	3662020	20.0
Triacontane, 1-br...	16.83	2.2	ng/ul	410274	6	15.69	3662020	20.0