

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
 Data File : BP020383.D
 Acq On : 15 May 2024 02:25
 Operator : MA/JU
 Sample : P2372-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43W0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020383.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	15	17	23	rVB4	10060	15165	0.88%	0.082%
2	3.487	64	68	71	rBV3	8583	13359	0.77%	0.072%
3	3.611	86	89	100	rBV	35461	80038	4.63%	0.432%
4	3.881	132	135	139	rVB4	7542	10186	0.59%	0.055%
5	4.187	185	187	194	rVB8	2011	2530	0.15%	0.014%
6	4.646	261	265	267	rBV5	3169	4034	0.23%	0.022%
7	4.758	282	284	292	rVB4	3513	6010	0.35%	0.032%
8	5.111	339	344	346	rBV6	2857	4370	0.25%	0.024%
9	5.428	396	398	402	rBV5	1639	2333	0.13%	0.013%
10	5.634	430	433	436	rBV4	2018	3293	0.19%	0.018%
11	5.887	473	476	481	rVB6	3134	5223	0.30%	0.028%
12	6.275	535	542	549	rVV3	13082	23240	1.34%	0.126%
13	6.699	611	614	617	rBV5	2408	3152	0.18%	0.017%
14	6.799	624	631	641	rBV2	88930	243026	14.05%	1.313%
15	6.887	641	646	652	rVV2	36827	71973	4.16%	0.389%
16	6.952	652	657	665	rVV2	97437	184791	10.69%	0.999%
17	7.011	665	667	673	rVV4	11390	17293	1.00%	0.093%
18	7.134	682	688	703	rBV	129084	257368	14.88%	1.391%
19	7.593	760	766	777	rBV	214736	387438	22.41%	2.094%
20	7.793	794	800	807	rBV4	19295	34480	1.99%	0.186%
21	7.840	807	808	814	rVB6	1689	2440	0.14%	0.013%
22	8.322	881	890	905	rBV2	50894	143716	8.31%	0.777%
23	8.440	905	910	930	rVB	53424	124355	7.19%	0.672%
24	8.764	958	965	982	rBV	132812	289435	16.74%	1.564%
25	9.093	1018	1021	1029	rVB9	3643	6254	0.36%	0.034%
26	9.263	1047	1050	1054	rBV5	1369	2329	0.13%	0.013%
27	9.369	1060	1068	1071	rBV7	4850	8712	0.50%	0.047%
28	9.475	1078	1086	1099	rBV	110409	240033	13.88%	1.297%
29	9.746	1123	1132	1133	rBV9	4182	8140	0.47%	0.044%
30	9.869	1149	1153	1157	rVB7	4567	6669	0.39%	0.036%
31	10.028	1171	1180	1205	rBV	109446	346217	20.02%	1.871%
32	10.228	1205	1214	1227	rVB8	6550	23169	1.34%	0.125%
33	10.369	1227	1238	1253	rVB	321623	587195	33.96%	3.173%
34	10.510	1257	1262	1263	rBV5	2043	2945	0.17%	0.016%
35	10.558	1263	1270	1287	rBV	53500	159478	9.22%	0.862%
36	10.905	1325	1329	1336	rBV10	4006	7987	0.46%	0.043%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
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37	10.987	1339	1343	1348	rBV8	4101	7864	0.45%	0.042%
38	11.175	1367	1375	1381	rVV6	13331	37322	2.16%	0.202%
39	11.375	1402	1409	1417	rVB8	4770	9666	0.56%	0.052%
40	11.593	1441	1446	1449	rBV7	2017	3878	0.22%	0.021%
41	11.793	1475	1480	1483	rVV4	5754	8898	0.51%	0.048%
42	11.987	1508	1513	1521	rVV10	4526	11198	0.65%	0.061%
43	12.052	1521	1524	1526	rVV3	2781	3977	0.23%	0.021%
44	12.069	1526	1527	1535	rVB8	3168	5288	0.31%	0.029%
45	12.569	1608	1612	1616	rBV7	3054	5355	0.31%	0.029%
46	12.710	1631	1636	1640	rBV7	2886	5062	0.29%	0.027%
47	12.763	1640	1645	1649	rVV4	5413	8040	0.46%	0.043%
48	13.057	1688	1695	1699	rBV3	14777	24383	1.41%	0.132%
49	13.099	1699	1702	1708	rVB8	4452	7540	0.44%	0.041%
50	13.575	1778	1783	1785	rBV6	4203	7209	0.42%	0.039%
51	13.640	1791	1794	1796	rBV4	4649	5945	0.34%	0.032%
52	13.687	1796	1802	1814	rVV	419851	711003	41.12%	3.842%
53	13.769	1814	1816	1823	rVB8	4883	8314	0.48%	0.045%
54	13.857	1828	1831	1834	rBV5	2904	3257	0.19%	0.018%
55	13.951	1838	1847	1858	rBV	471970	795022	45.98%	4.296%
56	14.163	1878	1883	1888	rBV2	25833	35505	2.05%	0.192%
57	14.204	1888	1890	1893	rBV4	3478	4750	0.27%	0.026%
58	14.263	1893	1900	1908	rBV	531790	876381	50.68%	4.736%
59	14.322	1908	1910	1913	rVB4	5989	5586	0.32%	0.030%
60	14.357	1913	1916	1923	rVB4	21961	34702	2.01%	0.188%
61	14.504	1932	1941	1946	rBV4	25807	63453	3.67%	0.343%
62	14.563	1946	1951	1963	rBV2	63202	232291	13.43%	1.255%
63	14.763	1982	1985	1986	rBV3	4545	5857	0.34%	0.032%
64	14.910	2006	2010	2015	rVB8	3977	6670	0.39%	0.036%
65	15.075	2031	2038	2043	rBV7	26670	50637	2.93%	0.274%
66	15.140	2045	2049	2056	rVB2	39984	56400	3.26%	0.305%
67	15.293	2068	2075	2083	rBV	692247	1074885	62.16%	5.808%
68	15.457	2097	2103	2114	rBV2	161888	322088	18.63%	1.740%
69	15.563	2116	2121	2130	rVB3	28363	62380	3.61%	0.337%
70	15.716	2143	2147	2158	rBV	90379	154237	8.92%	0.833%
71	15.869	2170	2173	2174	rBV3	4002	3592	0.21%	0.019%
72	16.040	2197	2202	2210	rBV3	67510	168579	9.75%	0.911%
73	16.163	2219	2223	2229	rBV9	4609	9674	0.56%	0.052%
74	16.516	2279	2283	2285	rBV5	9212	13085	0.76%	0.071%
75	16.645	2300	2305	2312	rBV8	16677	36758	2.13%	0.199%
76	16.728	2315	2319	2322	rVV4	17578	29671	1.72%	0.160%

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77	16.763	2322	2325	2335	rVV2	42132	93677	5.42%	0.506%
78	16.863	2335	2342	2346	rVV	70777	122317	7.07%	0.661%
79	16.916	2346	2351	2360	rVB2	42001	84800	4.90%	0.458%
80	17.104	2377	2383	2389	rBV	749597	1188700	68.75%	6.423%
81	17.210	2395	2401	2413	rVB	775837	1202339	69.54%	6.497%
82	17.310	2415	2418	2423	rVV7	12206	17824	1.03%	0.096%
83	17.545	2455	2458	2465	rBV7	17494	34456	1.99%	0.186%
84	17.640	2470	2474	2480	rVB2	71667	95269	5.51%	0.515%
85	17.798	2499	2501	2505	rBV4	15295	18689	1.08%	0.101%
86	17.928	2519	2523	2532	rBV4	28390	60261	3.49%	0.326%
87	18.004	2533	2536	2540	rBV5	20571	28096	1.62%	0.152%
88	18.069	2542	2547	2556	rBV2	212289	359675	20.80%	1.944%
89	18.369	2595	2598	2603	rBV	79164	91620	5.30%	0.495%
90	19.039	2709	2712	2716	rVB2	74391	91818	5.31%	0.496%
91	19.087	2717	2720	2726	rBV3	85216	112947	6.53%	0.610%
92	19.163	2729	2733	2736	rBV6	36990	61697	3.57%	0.333%
93	19.416	2773	2776	2785	rVB3	86893	142116	8.22%	0.768%
94	19.610	2804	2809	2820	rBV	1094638	1729113	100.00%	9.343%
95	20.222	2909	2913	2917	rBV4	77686	122457	7.08%	0.662%
96	20.875	3022	3024	3032	rVB7	85110	127146	7.35%	0.687%
97	21.286	3091	3094	3100	rVB2	174017	262735	15.19%	1.420%
98	21.604	3143	3148	3157	rVB	822138	1333314	77.11%	7.205%
99	24.733	3672	3680	3687	rBV	603486	1491118	86.24%	8.057%
100	24.951	3711	3717	3729	rVB	490808	1389388	80.35%	7.508%

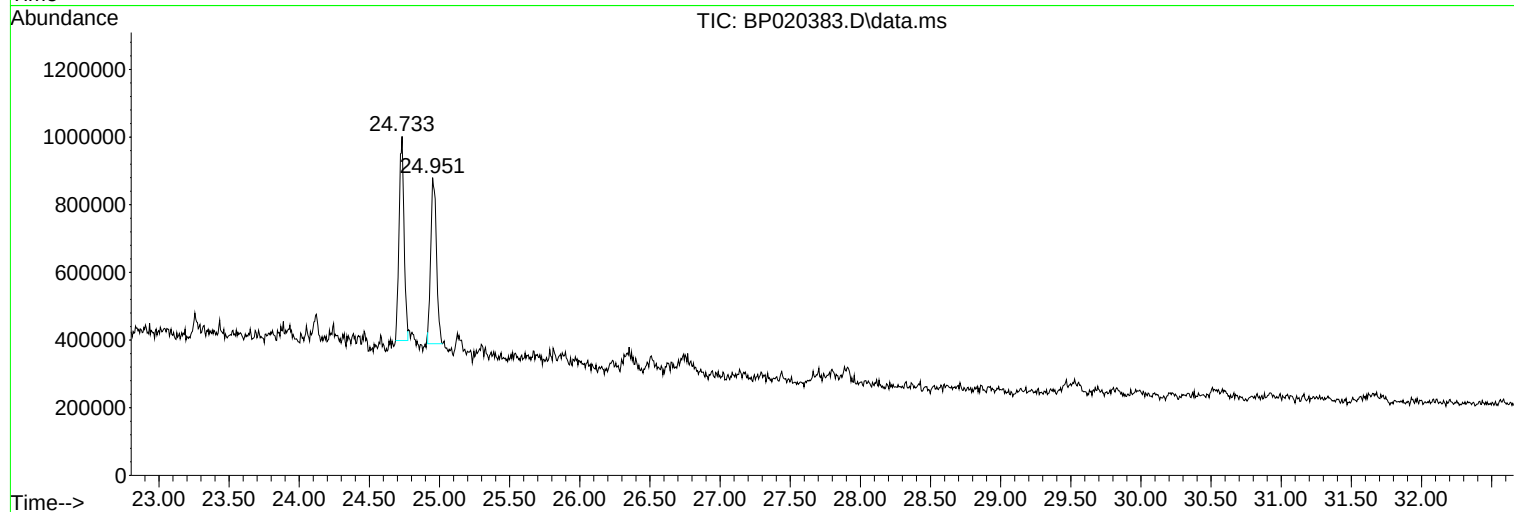
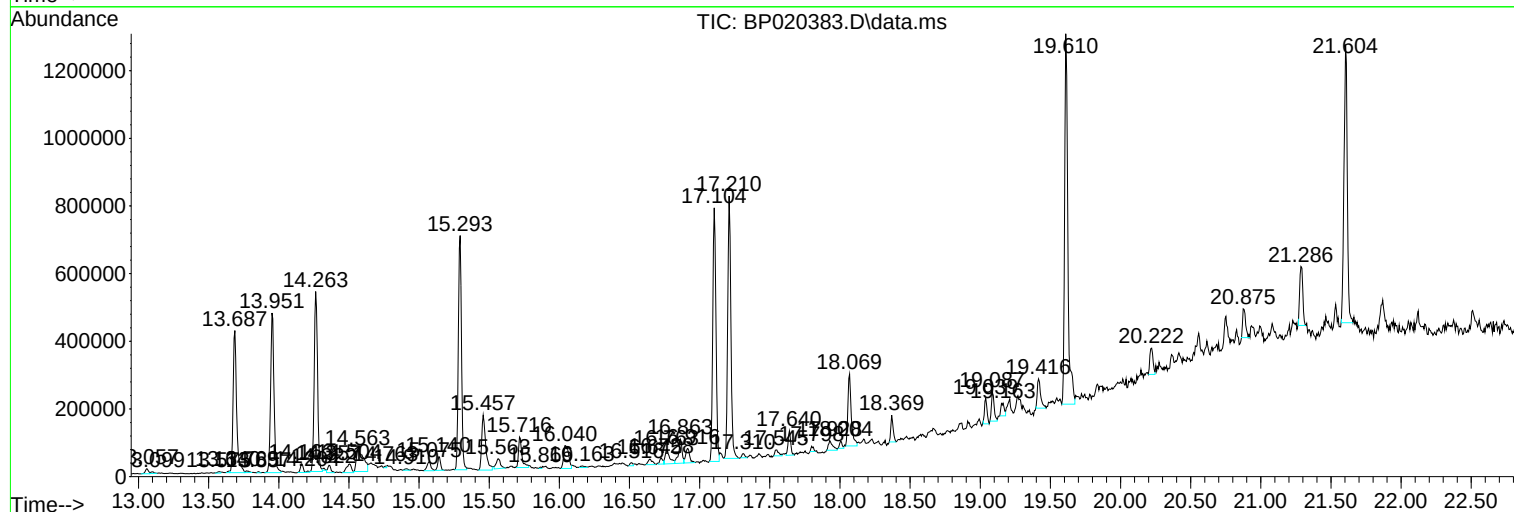
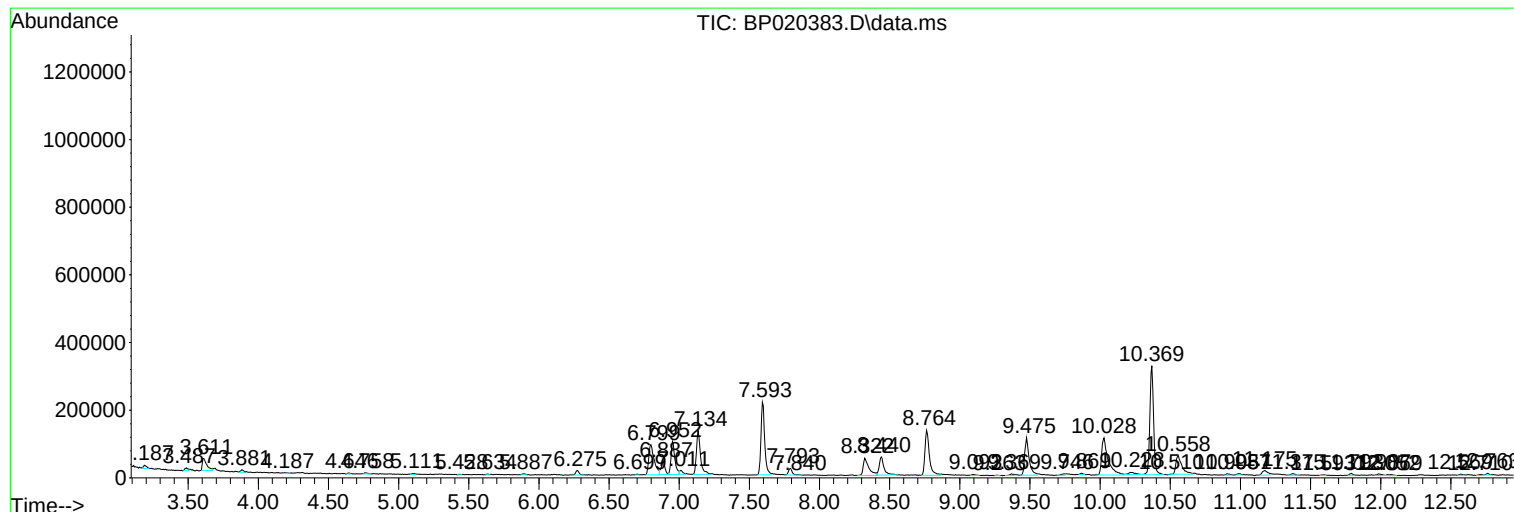
Sum of corrected areas: 18506350

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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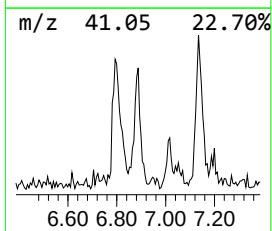
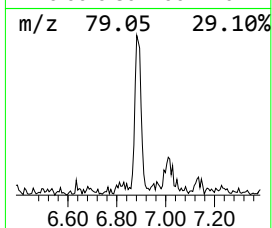
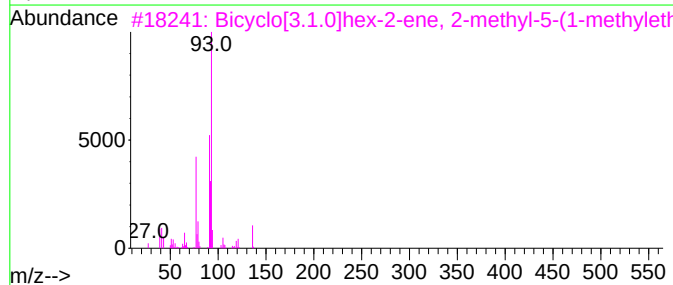
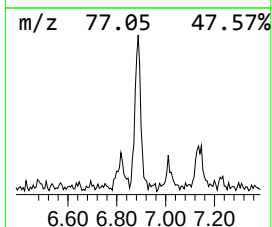
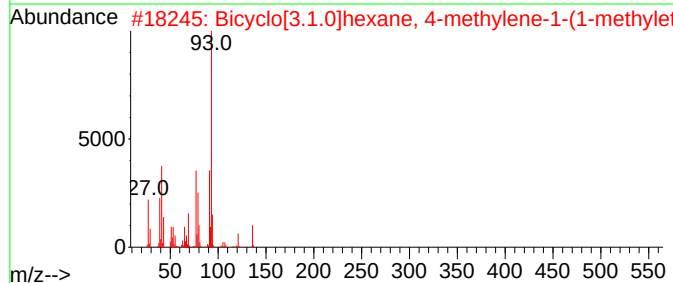
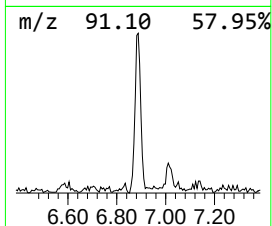
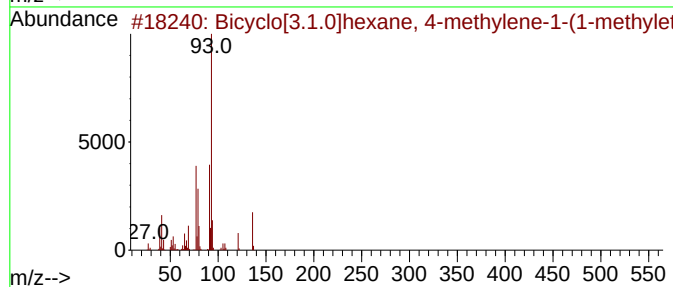
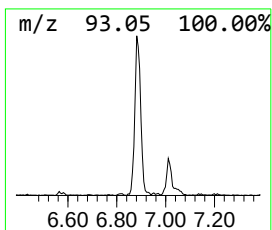
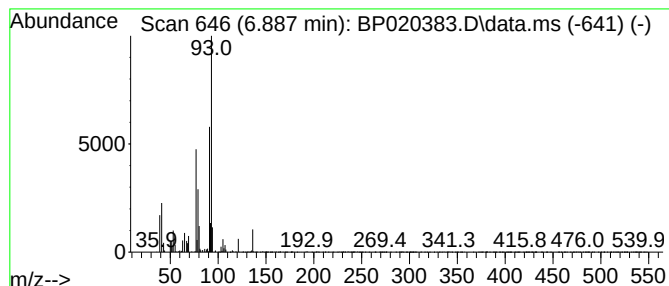
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Bicyclo[3.1.0]hexane, 4-met... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.887	3.72 ng/ul	71973	1,4-Dichlorobenzene-d4	7.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	91
2			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	90
3			Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	87
4			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	83
5			(1S)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-26-4	80



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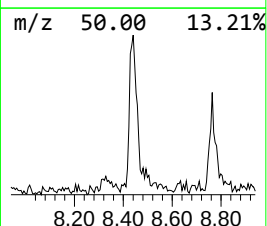
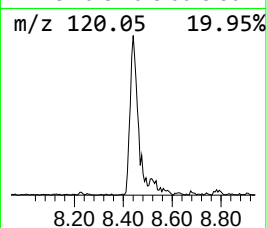
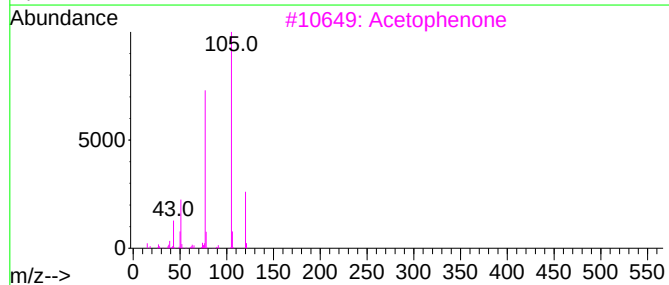
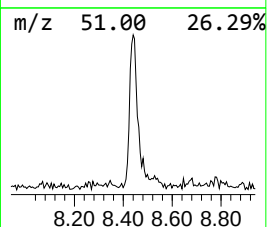
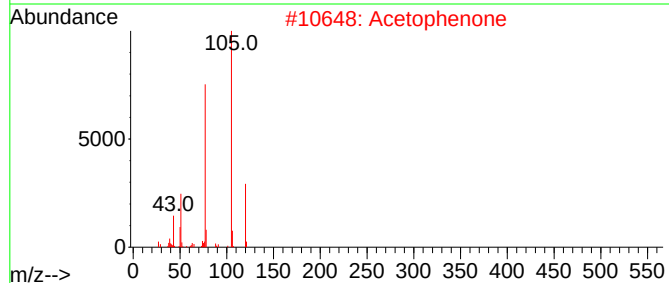
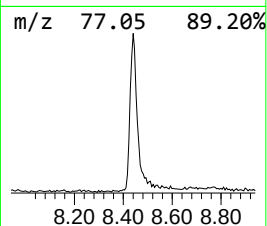
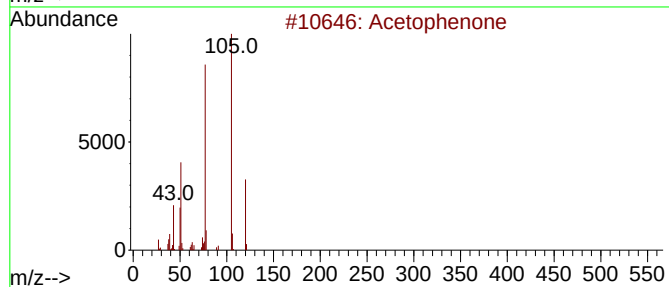
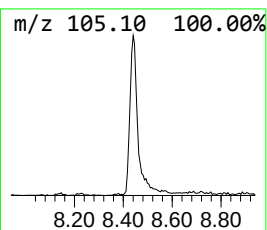
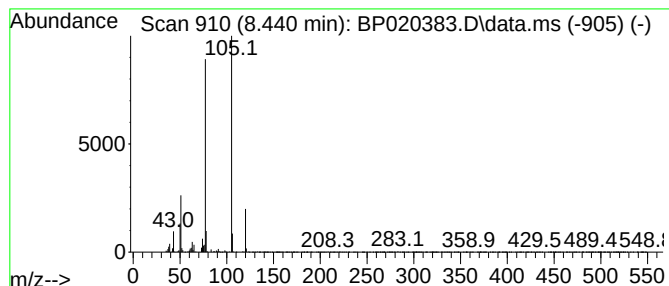
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Aceto phenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.440	6.42 ng/ul	124355	1,4-Dichlorobenzene-d4	7.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	91
2			Acetophenone	120	C8H8O	000098-86-2	91
3			Acetophenone	120	C8H8O	000098-86-2	91
4			Acetophenone	120	C8H8O	000098-86-2	91
5			1-Butanone, 3-methyl-1-phenyl-	162	C11H14O	000582-62-7	78



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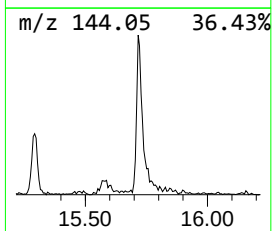
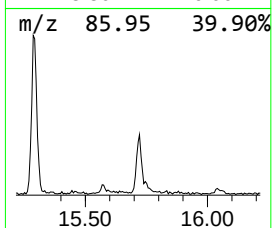
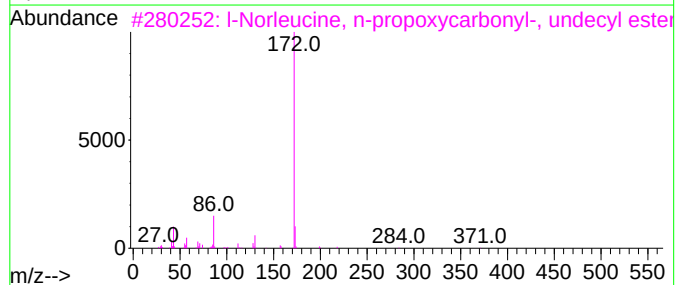
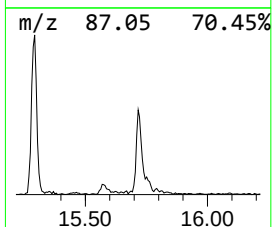
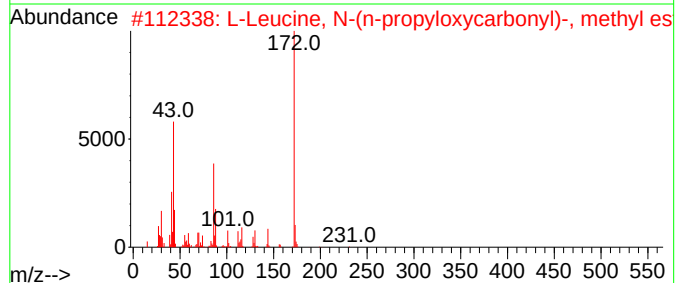
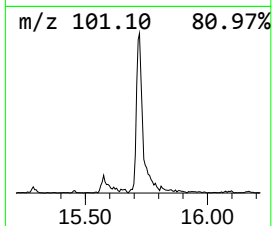
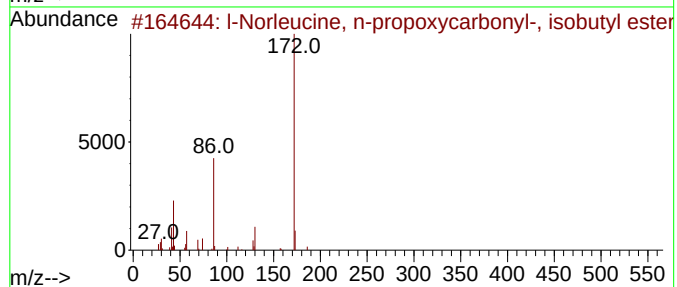
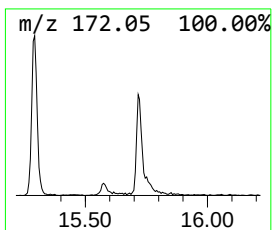
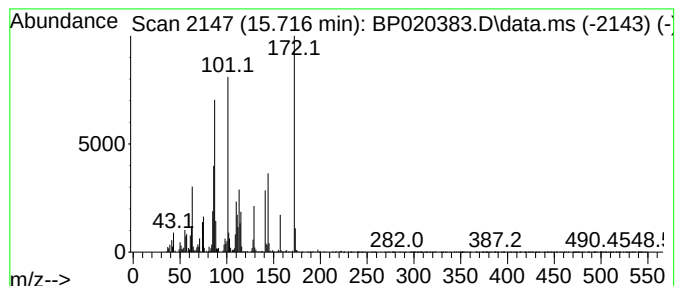
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.716	2.60 ng/ul	154237	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			l-Norleucine, n-propoxycarbonyl-...	273	C14H27NO4	1000329-53-8	22
2			L-Leucine, N-(n-propyloxycarbonyl-...	231	C11H21NO4	1000452-30-4	22
3			l-Norleucine, n-propoxycarbonyl-...	371	C21H41NO4	1000329-54-4	14
4			l-Norleucine, n-propoxycarbonyl-...	385	C22H43NO4	1000329-54-5	14
5			Carbamic acid, N-(2-butyl)-N-pro...	201	C11H23NO2	1000491-05-7	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

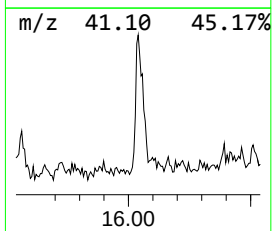
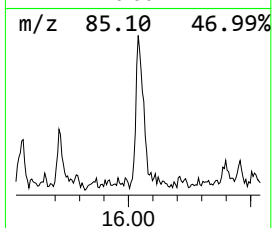
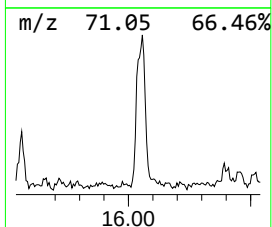
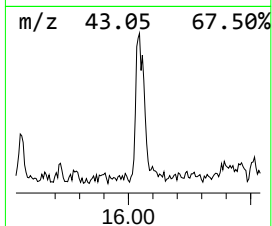
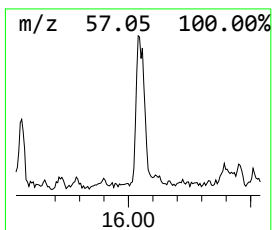
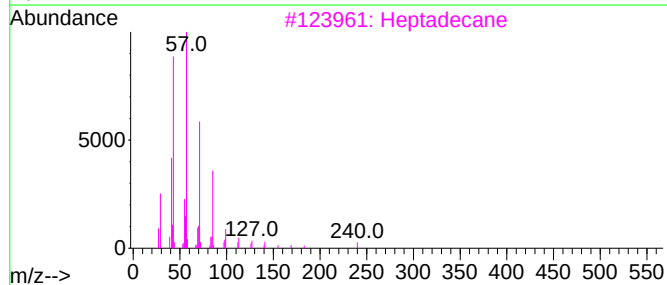
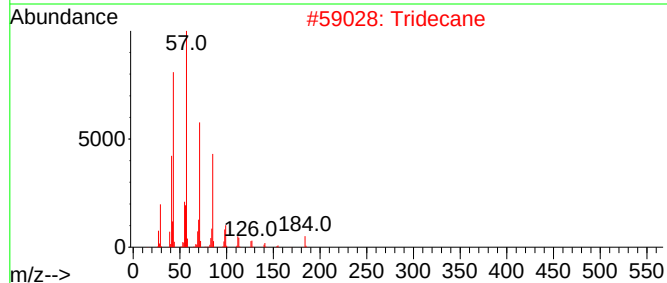
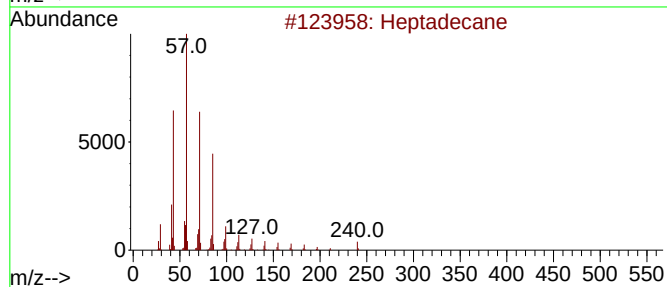
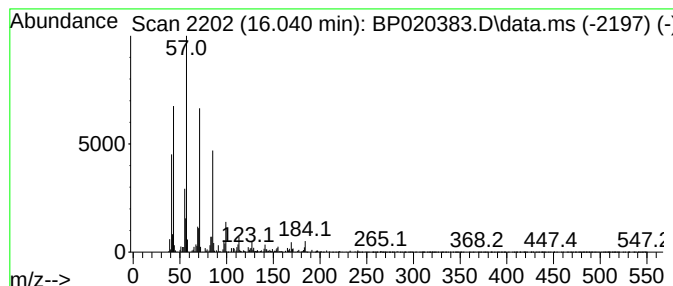
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.040	2.84 ng/ul	168579	Phenanthrene-d10		17.104	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	94
2	Tridecane		184	C13H28	000629-50-5	93
3	Heptadecane		240	C17H36	000629-78-7	93
4	Heptadecane		240	C17H36	000629-78-7	91
5	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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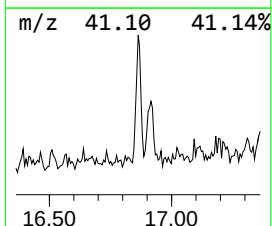
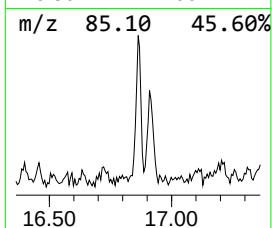
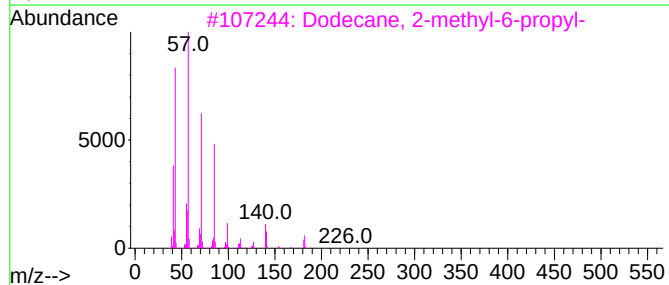
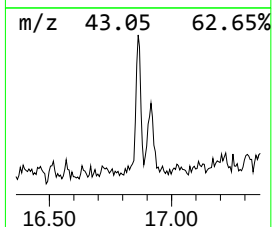
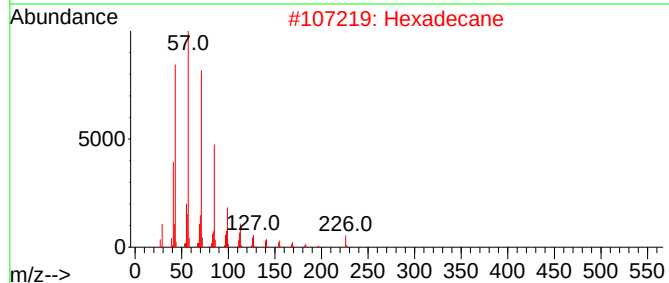
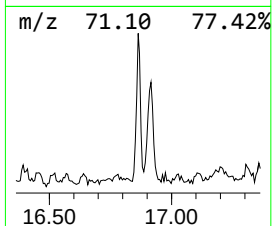
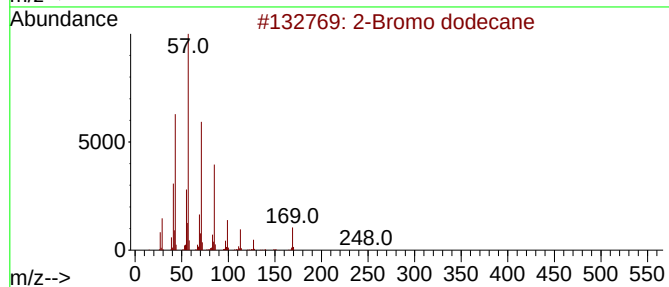
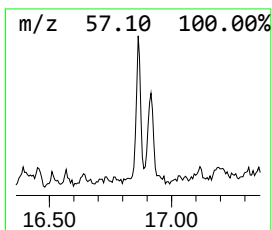
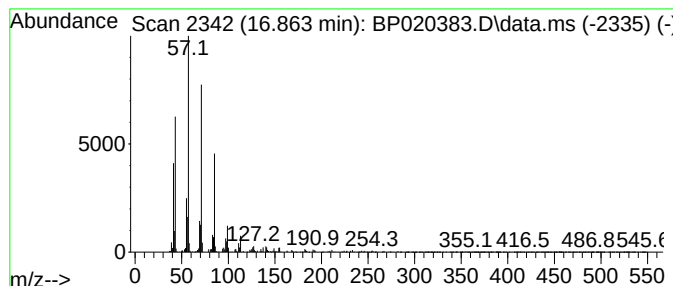
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Bromo dodecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.863	2.06 ng/ul	122317	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Bromo dodecane	248	C12H25Br	013187-99-0	94
2			Hexadecane	226	C16H34	000544-76-3	94
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
4			Eicosane	282	C20H42	000112-95-8	91
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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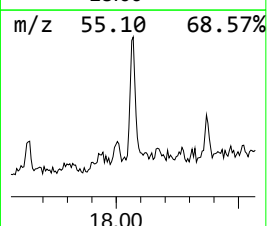
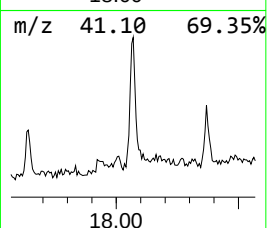
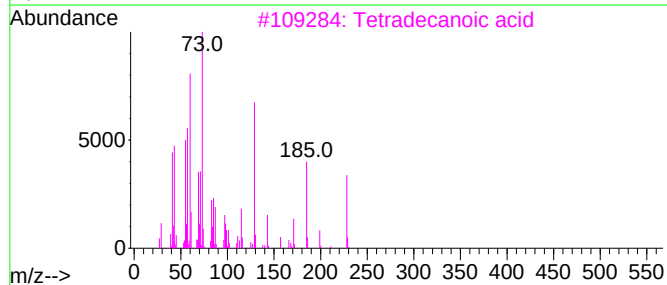
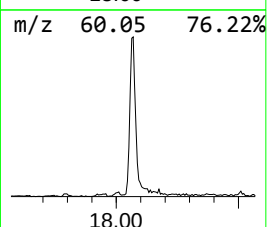
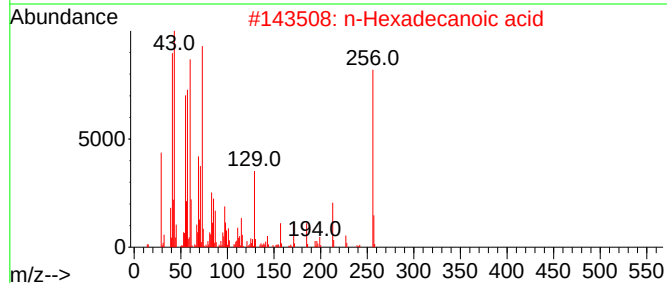
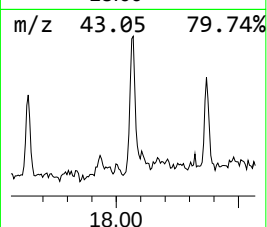
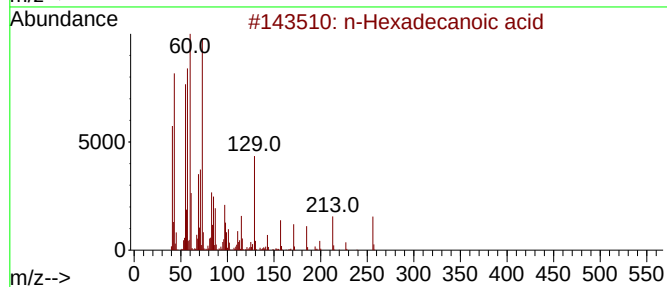
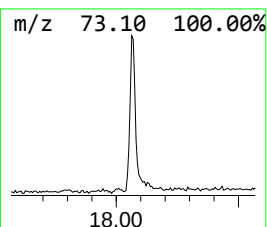
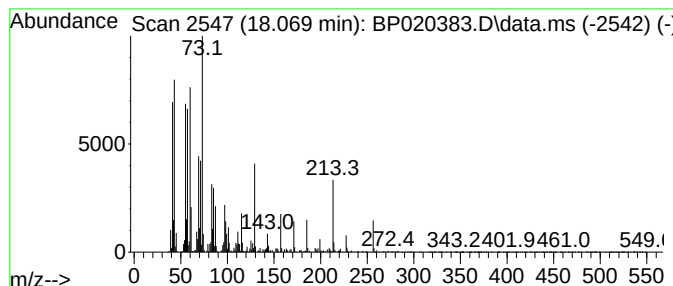
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.069	6.05 ng/ul	359675	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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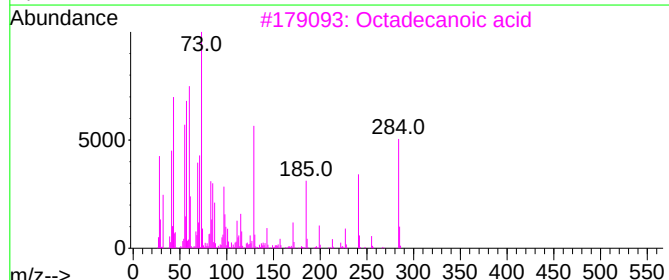
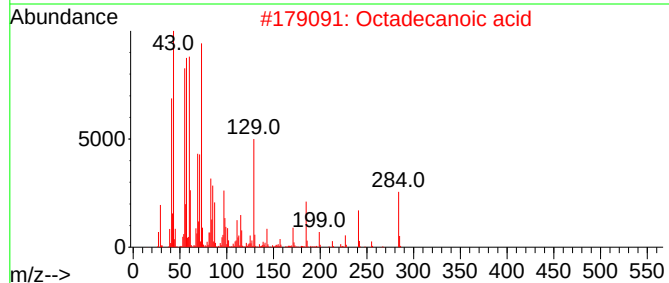
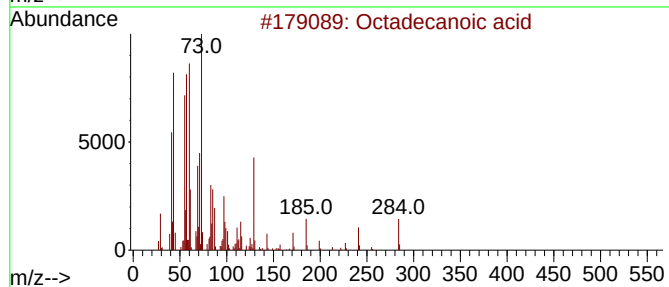
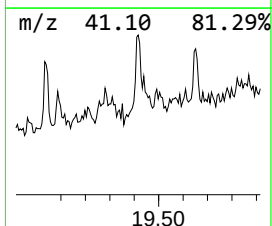
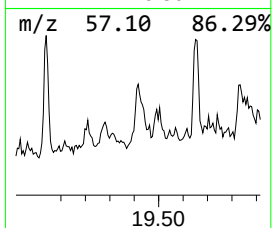
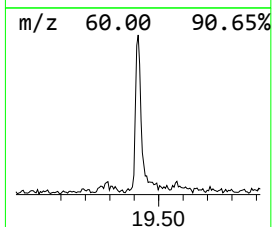
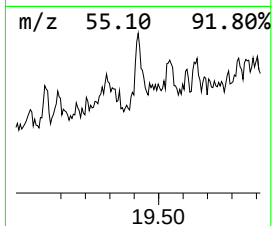
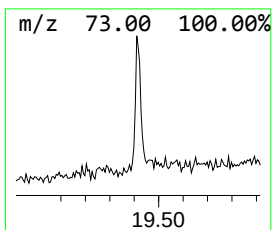
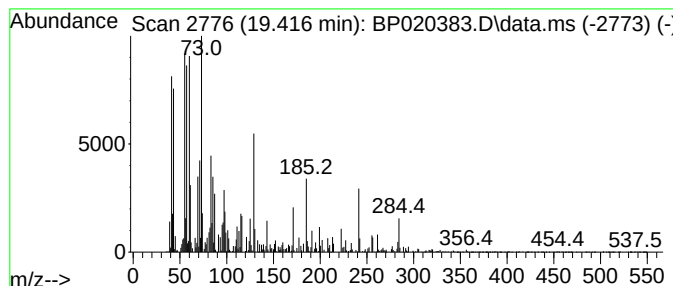
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.416	2.13 ng/ul	142116	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	97
2			Octadecanoic acid	284	C18H36O2	000057-11-4	96
3			Octadecanoic acid	284	C18H36O2	000057-11-4	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

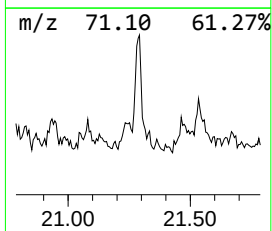
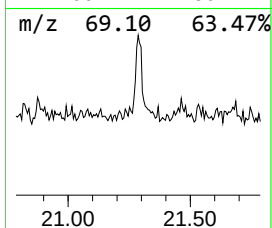
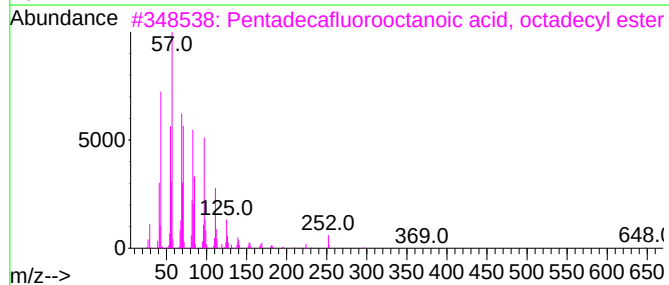
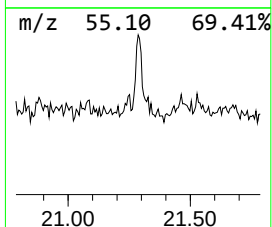
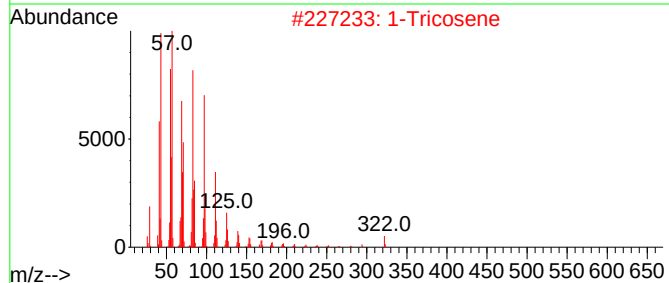
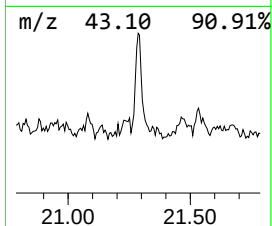
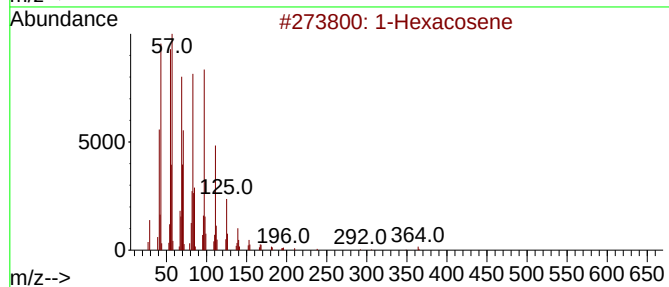
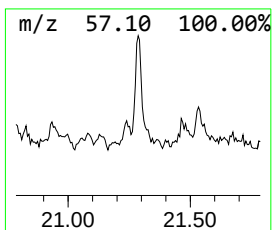
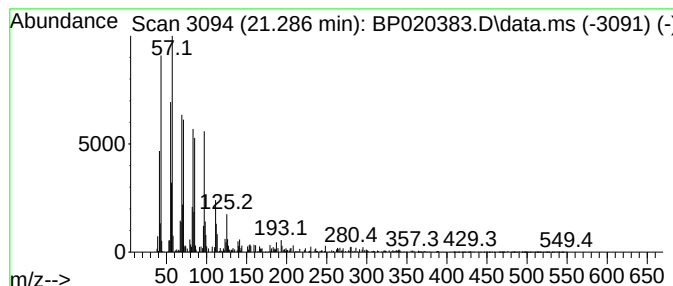
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.286	3.94 ng/ul	262735	Chrysene-d12		21.604	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Hexacosene		364	C26H52	018835-33-1	97
2	1-Tricosene		322	C23H46	018835-32-0	96
3	Pentadecafluorooctanoic acid, oc...		666	C26H37F15O2	1000406-04-8	90
4	1-Docosene		308	C22H44	001599-67-3	89
5	Z-5-Nonadecene		266	C19H38	1000131-11-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020383.D
Acq On : 15 May 2024 02:25
Operator : MA/JU
Sample : P2372-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Aceto phenone	8.440	6.4	ng/ul	124355	1	7.593	387438	20.0
unknown-01	15.716	2.6	ng/ul	154237	4	17.104	1188700	20.0
(DEL) Alkane: S...	16.040	2.8	ng/ul	168579	4	17.104	1188700	20.0
2-Bromo dodecane	16.863	2.1	ng/ul	122317	4	17.104	1188700	20.0
n-Hexadecanoic ...	18.069	6.0	ng/ul	359675	4	17.104	1188700	20.0
Octadecanoic acid	19.416	2.1	ng/ul	142116	5	21.604	1333310	20.0
(DEL) Alkane: S...	21.286	3.9	ng/ul	262735	5	21.604	1333310	20.0