

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006566.D
 Acq On : 07 Aug 2021 13:50
 Operator : CG/JU
 Sample : M3163-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGEX4

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006566.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.681	98	101	111	rBV	266351	450529	17.18%	1.148%
2	6.928	647	653	663	rVB	520720	849916	32.42%	2.166%
3	7.099	676	682	692	rVB	429976	700320	26.71%	1.785%
4	7.287	708	714	721	rVB	540034	851806	32.49%	2.171%
5	7.752	787	793	801	rVB	698987	1094433	41.74%	2.789%
6	8.457	907	913	921	rBV	572634	900659	34.35%	2.295%
7	8.575	930	933	940	rVB2	13018	21980	0.84%	0.056%
8	8.904	983	989	999	rVB	527611	853237	32.54%	2.174%
9	9.328	1058	1061	1068	rVB2	14515	22435	0.86%	0.057%
10	9.622	1105	1111	1120	rVB	406517	651688	24.86%	1.661%
11	10.157	1195	1202	1221	rBV	550501	927322	35.37%	2.363%
12	10.328	1225	1231	1243	rVV	170685	302179	11.53%	0.770%
13	10.534	1259	1266	1274	rVB	936603	1553566	59.25%	3.959%
14	10.675	1282	1290	1302	rBV	467829	843519	32.17%	2.150%
15	11.010	1343	1347	1351	rBV6	3522	5981	0.23%	0.015%
16	11.563	1437	1441	1445	rVB7	2509	4142	0.16%	0.011%
17	11.816	1475	1484	1490	rBV4	9396	22312	0.85%	0.057%
18	11.940	1503	1505	1508	rBV4	2020	2901	0.11%	0.007%
19	12.045	1519	1523	1524	rBV3	4650	5109	0.19%	0.013%
20	12.128	1532	1537	1541	rVB4	11985	19609	0.75%	0.050%
21	12.745	1637	1642	1645	rBV6	2767	5530	0.21%	0.014%
22	12.787	1647	1649	1655	rVB7	2553	3616	0.14%	0.009%
23	12.857	1658	1661	1664	rBV4	2838	4013	0.15%	0.010%
24	12.992	1681	1684	1687	rBV5	5370	6087	0.23%	0.016%
25	13.104	1698	1703	1707	rBV8	3510	6039	0.23%	0.015%
26	13.163	1709	1713	1715	rVV5	2275	2711	0.10%	0.007%
27	13.216	1718	1722	1726	rBV2	9135	12433	0.47%	0.032%
28	13.351	1741	1745	1751	rVB9	5732	10314	0.39%	0.026%
29	13.704	1799	1805	1814	rBV5	25057	39671	1.51%	0.101%
30	13.798	1814	1821	1831	rBV	1009469	1430271	54.55%	3.645%
31	13.875	1831	1834	1838	rVV5	8177	14690	0.56%	0.037%
32	13.910	1838	1840	1844	rVB5	3222	4662	0.18%	0.012%
33	14.069	1859	1867	1877	rBV	1109167	1640758	62.58%	4.181%
34	14.222	1889	1893	1894	rBV4	4390	5315	0.20%	0.014%
35	14.287	1902	1904	1909	rVB6	4848	5915	0.23%	0.015%
36	14.381	1909	1920	1928	rBV	1301603	1924980	73.42%	4.905%

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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 >
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37	14.445	1928	1931	1934	rVB5	3197	4670	0.18%	0.012%
38	14.551	1946	1949	1950	rBV3	3268	3550	0.14%	0.009%
39	14.586	1950	1955	1975	rVV	313790	594390	22.67%	1.515%
40	14.745	1979	1982	1989	rVV7	8378	14641	0.56%	0.037%
41	14.810	1989	1993	1999	rVB6	9937	12720	0.49%	0.032%
42	14.992	2022	2024	2029	rVB6	2502	3717	0.14%	0.009%
43	15.039	2029	2032	2036	rVB6	2124	3279	0.13%	0.008%
44	15.192	2054	2058	2064	rBV7	4146	7645	0.29%	0.019%
45	15.251	2064	2068	2072	rVB7	2535	3060	0.12%	0.008%
46	15.375	2082	2089	2096	rBV	1385602	1975846	75.36%	5.035%
47	15.492	2104	2109	2119	rBV	316993	459825	17.54%	1.172%
48	15.610	2126	2129	2134	rVB	17099	23626	0.90%	0.060%
49	15.763	2150	2155	2156	rBV5	2084	3480	0.13%	0.009%
50	15.833	2163	2167	2174	rVB9	7201	13640	0.52%	0.035%
51	15.945	2180	2186	2188	rBV7	5550	8066	0.31%	0.021%
52	16.110	2212	2214	2219	rVV6	2905	4677	0.18%	0.012%
53	16.157	2219	2222	2227	rVV7	3415	6096	0.23%	0.016%
54	16.281	2238	2243	2245	rBV6	4479	7639	0.29%	0.019%
55	16.345	2249	2254	2256	rVB6	3060	4288	0.16%	0.011%
56	16.433	2267	2269	2275	rVB7	2894	3152	0.12%	0.008%
57	16.486	2275	2278	2280	rVB4	3095	2850	0.11%	0.007%
58	16.516	2280	2283	2284	rBV3	1928	2692	0.10%	0.007%
59	16.545	2284	2288	2290	rBV4	3736	3771	0.14%	0.010%
60	16.633	2300	2303	2305	rBV3	2805	3113	0.12%	0.008%
61	16.739	2319	2321	2325	rBV5	3166	3864	0.15%	0.010%
62	16.775	2325	2327	2331	rBV5	2037	3166	0.12%	0.008%
63	16.910	2344	2350	2355	rBV6	9054	20430	0.78%	0.052%
64	17.133	2381	2388	2398	rBV	1519454	2173732	82.91%	5.539%
65	17.228	2398	2404	2417	rVV2	1400937	2043003	77.92%	5.206%
66	17.504	2449	2451	2456	rBV6	2914	5017	0.19%	0.013%
67	17.651	2473	2476	2477	rBV3	3035	2943	0.11%	0.007%
68	17.822	2501	2505	2514	rVB	28816	37746	1.44%	0.096%
69	18.039	2537	2542	2549	rBV	161348	249623	9.52%	0.636%
70	19.051	2710	2714	2718	rBV4	20603	29586	1.13%	0.075%
71	19.122	2722	2726	2729	rBV2	20859	30808	1.18%	0.079%
72	19.274	2748	2752	2758	rBV2	70201	102067	3.89%	0.260%
73	19.474	2780	2786	2793	rBV	1732539	2331954	88.94%	5.943%
74	20.022	2876	2879	2882	rVB2	21028	22855	0.87%	0.058%
75	20.657	2983	2987	2991	rBV4	80434	107657	4.11%	0.274%
76	20.727	2997	2999	3004	rVB6	35762	38296	1.46%	0.098%

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77	20.845	3015	3019	3023	rBV	331883	432340	16.49%	1.102%
78	20.886	3024	3026	3032	rVB6	69960	91999	3.51%	0.234%
79	20.980	3034	3042	3048	rBV2	932703	1548155	59.05%	3.945%
80	21.092	3057	3061	3066	rBV3	111443	159924	6.10%	0.408%
81	21.145	3066	3070	3074	rBV	606827	696824	26.58%	1.776%
82	21.192	3075	3078	3080	rVV2	179358	210563	8.03%	0.537%
83	21.227	3080	3084	3093	rVB	1863685	2409314	91.89%	6.140%
84	21.845	3185	3189	3201	rBV2	326755	699144	26.67%	1.782%
85	22.868	3358	3363	3368	rBV	707634	1045264	39.87%	2.664%
86	22.921	3368	3372	3383	rVB2	268606	548917	20.94%	1.399%
87	23.404	3448	3454	3461	rBV	1106573	2064682	78.75%	5.261%
88	23.557	3473	3480	3494	rVB2	1276851	2621939	100.00%	6.682%
89	24.180	3580	3586	3594	rVB	578464	1138545	43.42%	2.901%

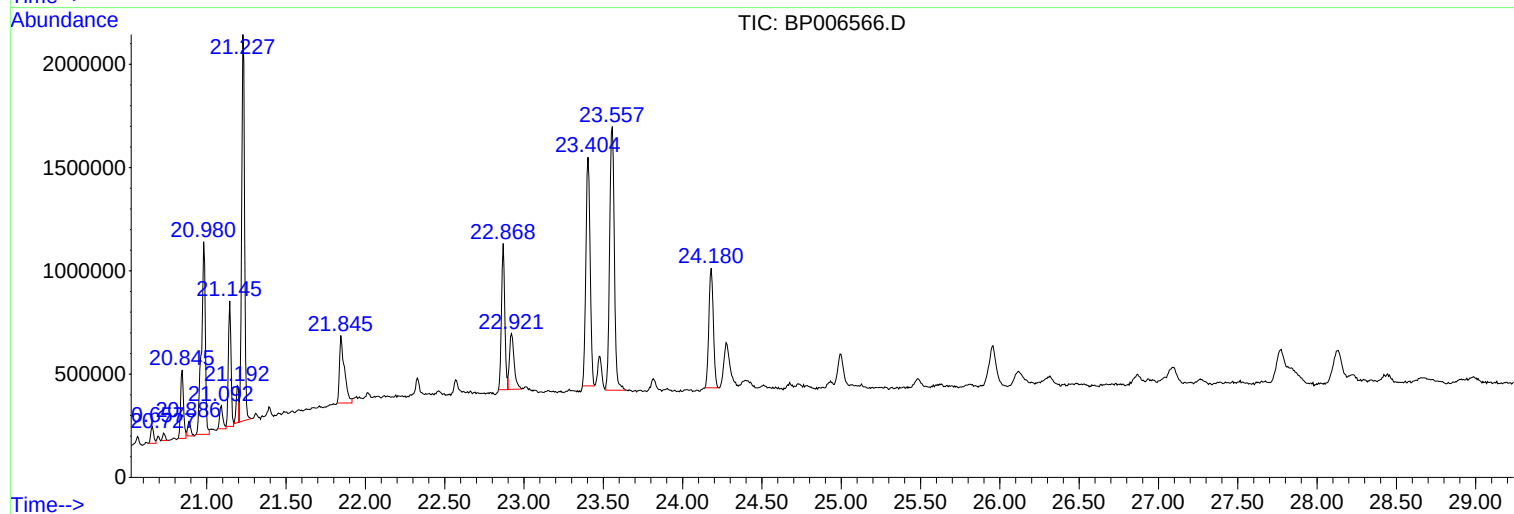
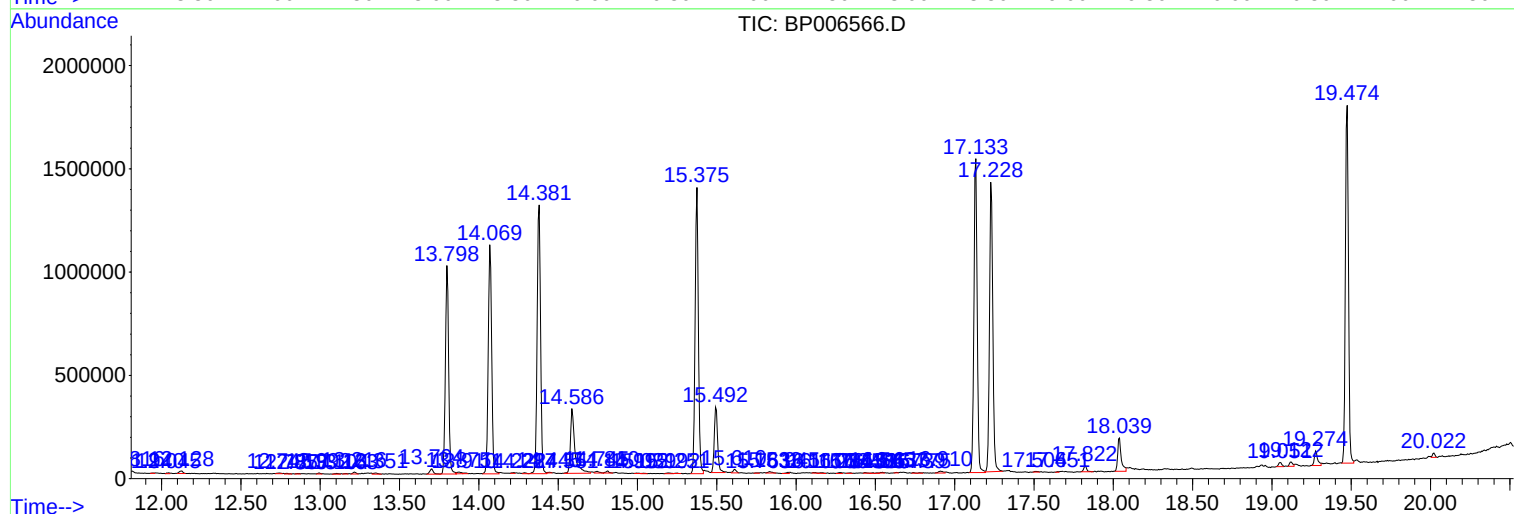
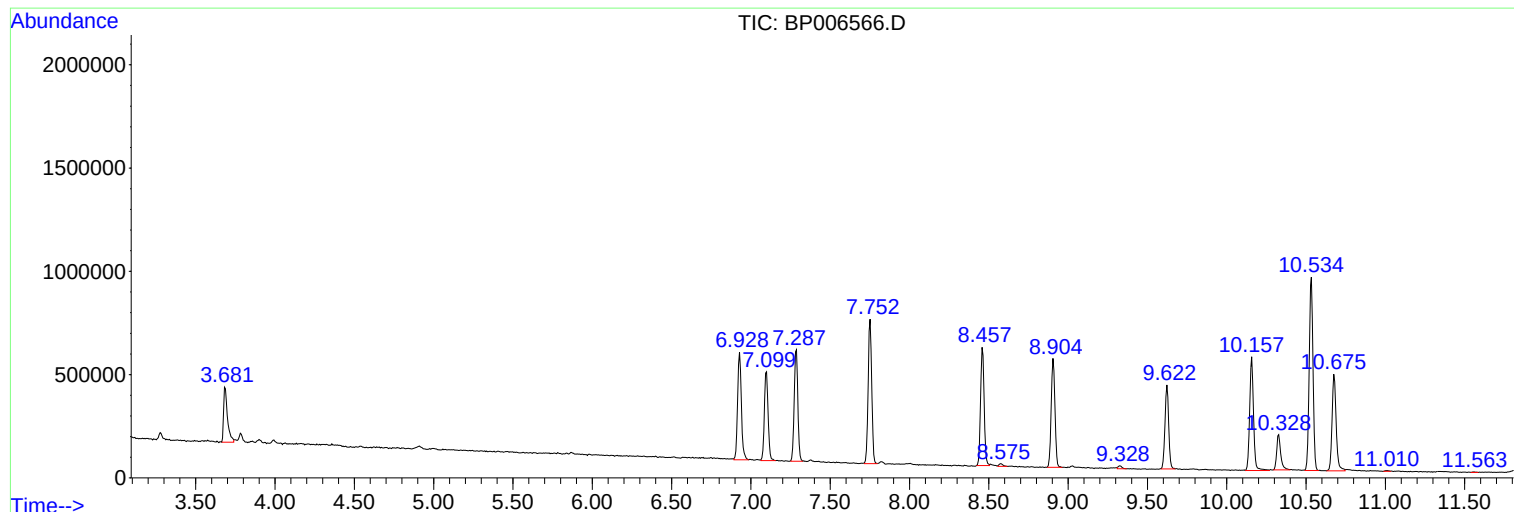
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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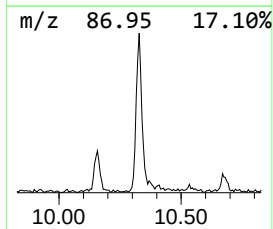
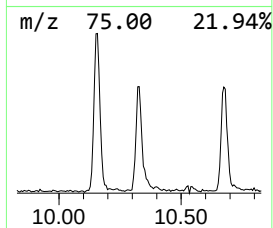
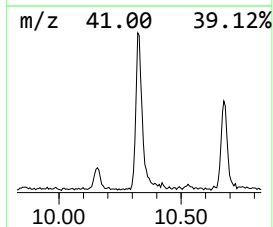
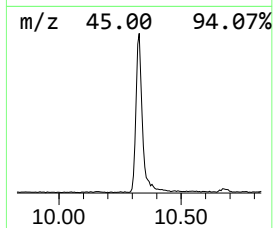
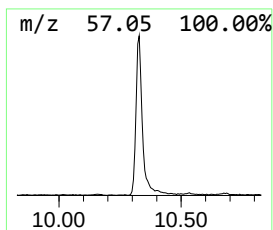
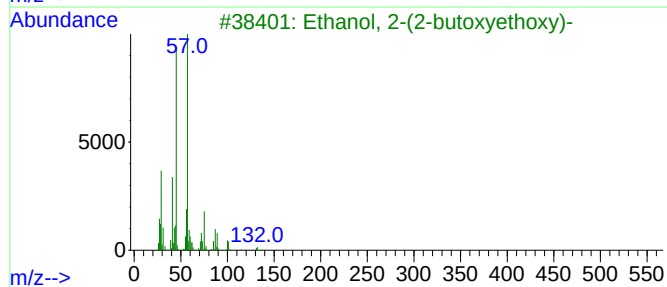
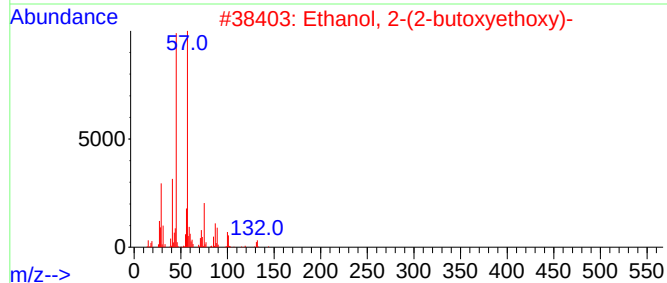
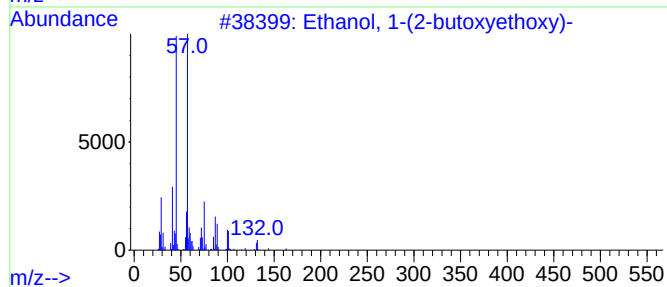
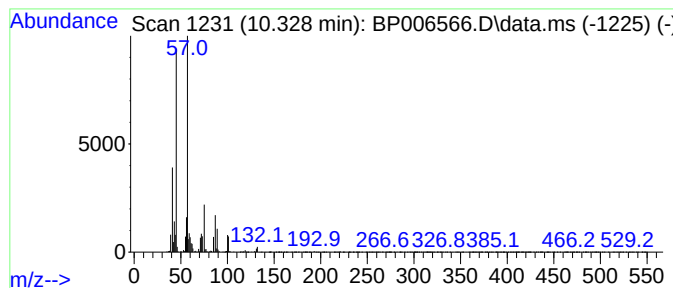
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.330	3.89 ng/ul	302179	Naphthalene-d8	10.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83



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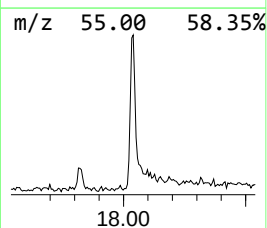
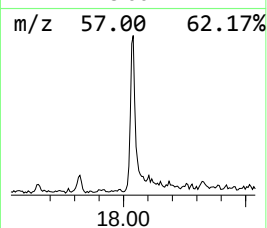
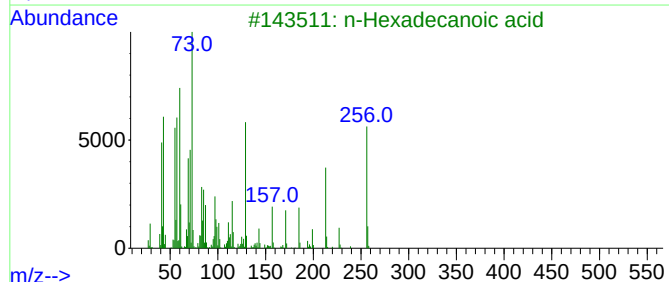
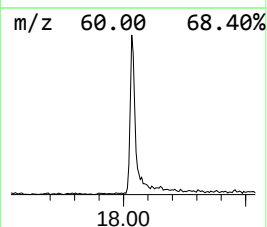
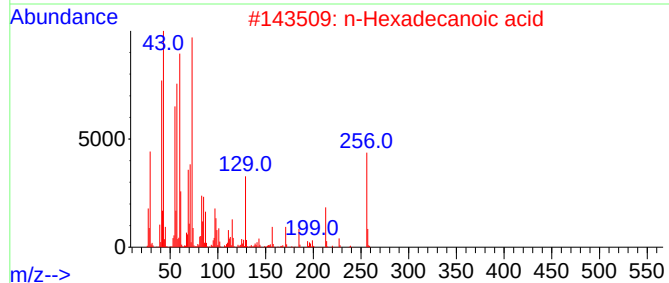
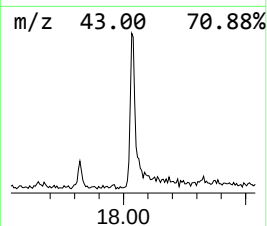
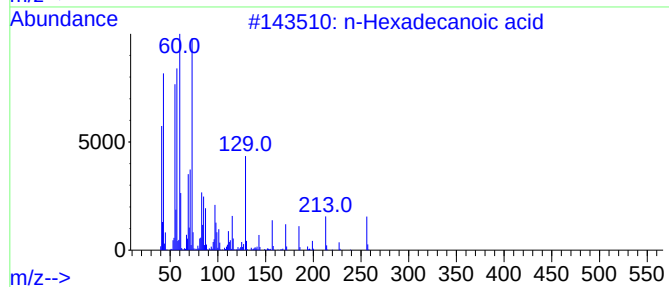
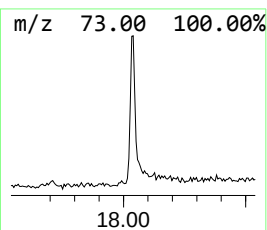
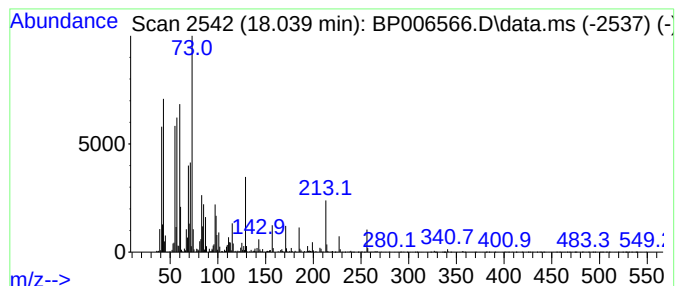
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.040	2.30 ng/ul	249623	Phenanthrene-d10	17.133

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
5			Octadecanoic acid	284	C18H36O2	000057-11-4	81



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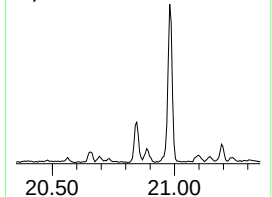
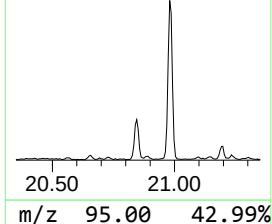
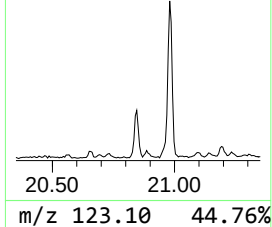
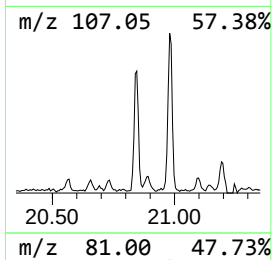
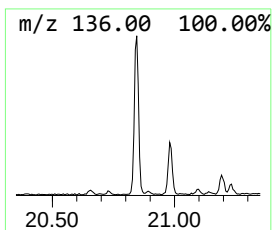
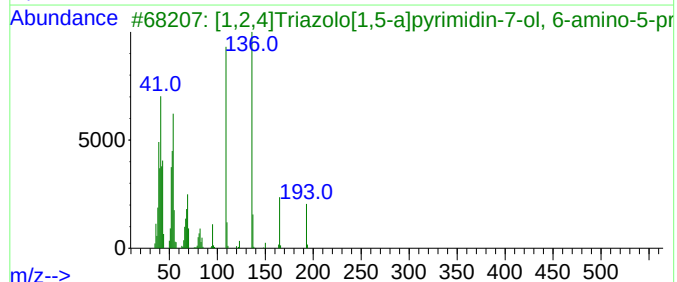
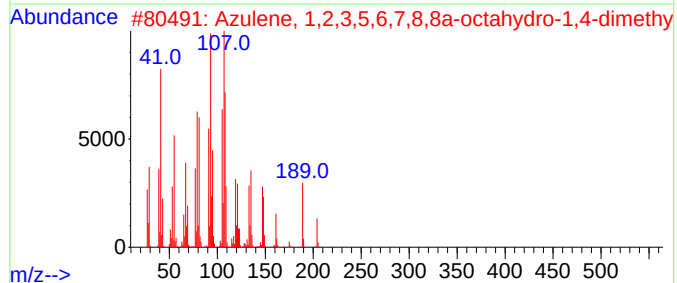
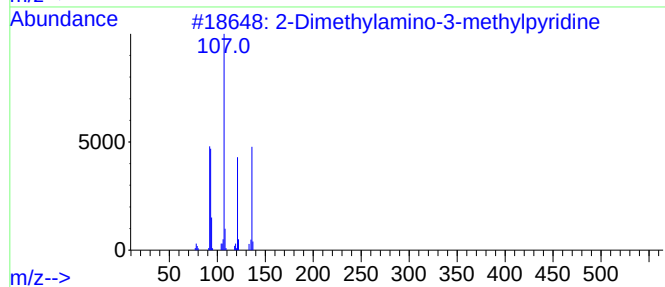
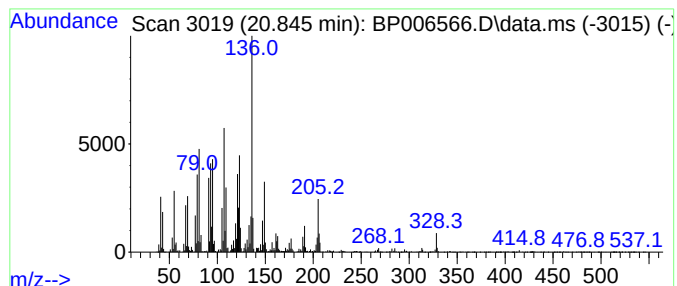
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Dimethylamino-3-methylpyr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.850	3.59 ng/ul	432340	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Dimethylamino-3-methylpyridine	136	C8H12N2	061713-46-0	50
2			Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	46
3			[1,2,4]Triazolo[1,5-a]pyrimidin-...	193	C8H11N5O	1000362-66-7	35
4			11,11-Dimethyl-4,8-dimethylenebi...	220	C15H24O	079580-01-1	30
5			Spiro[5.5]undeca-1,8-diene, 1,5,...	204	C15H24	019912-83-5	30



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ClientSampleId :
BGEX4

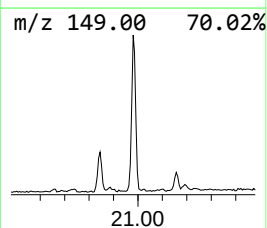
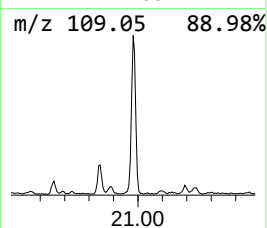
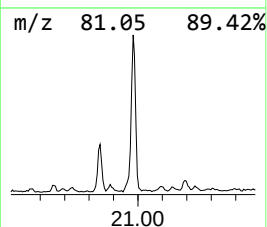
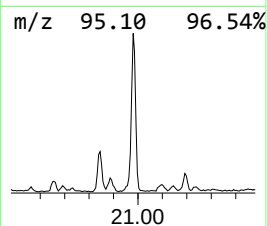
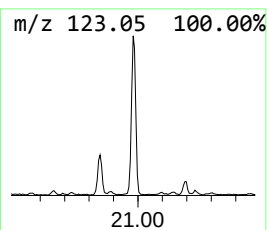
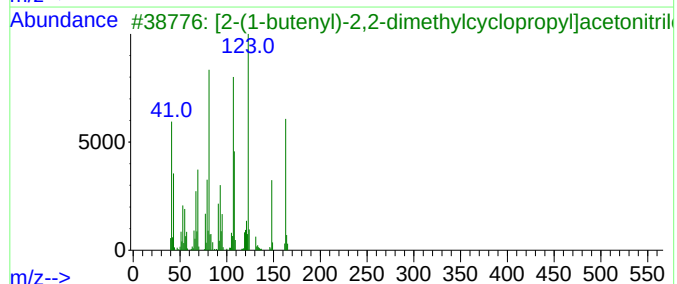
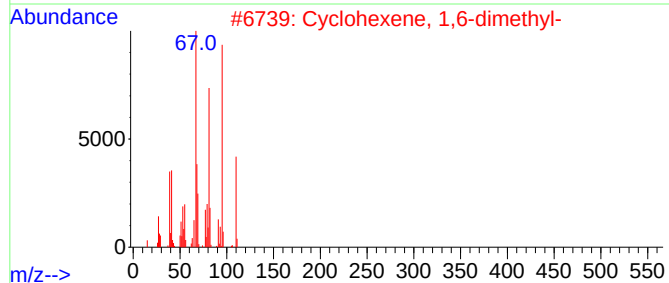
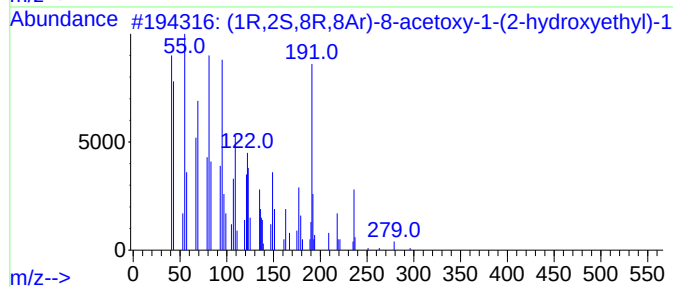
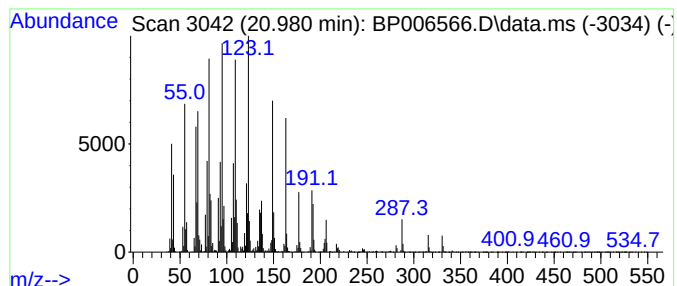
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (1R,2S,8R,8Ar)-8-acetoxy-1-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.980	12.85 ng/ul	1548160	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R,2S,8R,8Ar)-8-acetoxy-1-(2-hy...	296	C18H32O3	1000298-98-4	52		
2	Cyclohexene, 1,6-dimethyl-	110	C8H14	001759-64-4	35		
3	[2-(1-butenyl)-2,2-dimethylcyclo...	163	C11H17N	1000111-15-4	30		
4	Benzamide, 4-fluoro-N-(7-methyl-...	287	C14H10FN3OS	1000350-88-7	30		
5	(-)-Isolongifolol, chlorodifluor...	334	C17H25ClF2O2	1000447-45-0	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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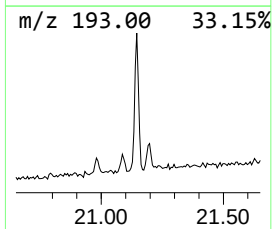
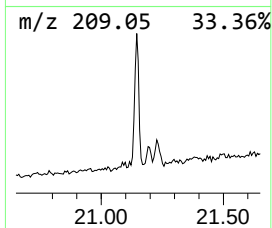
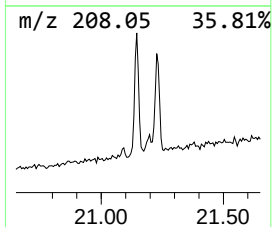
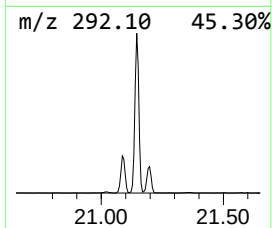
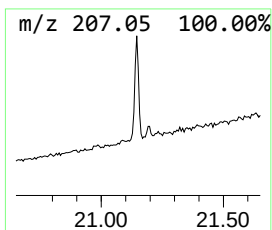
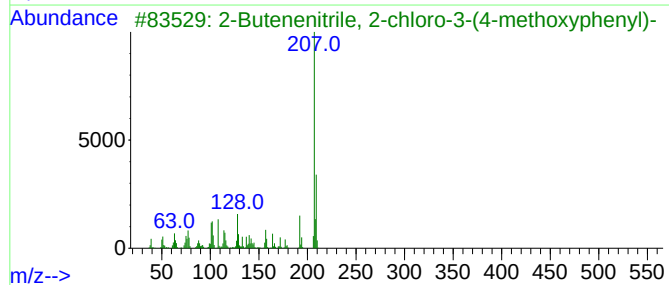
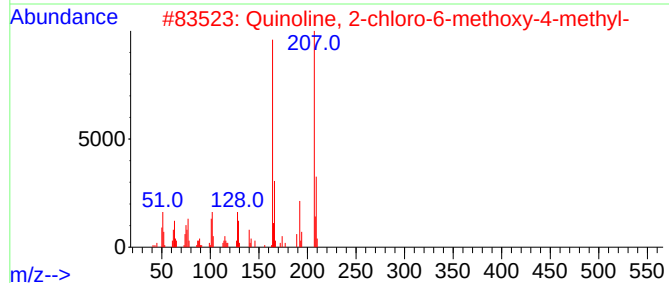
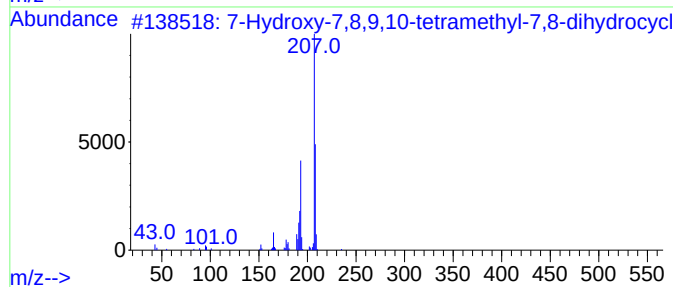
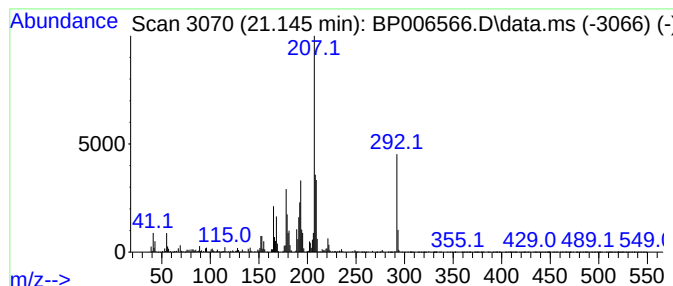
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.150	5.78 ng/ul	696824	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Hydroxy-7,8,9,10-tetramethyl-7...	252	C18H20O	1000110-34-9	47
2			Quinoline, 2-chloro-6-methoxy-4-...	207	C11H10ClNO	006340-55-2	43
3			2-Butenenitrile, 2-chloro-3-(4-m...	207	C11H10ClNO	1010305-66-7	38
4			Silicic acid, diethyl bis(trimet...	296	C10H28O4Si3	003555-45-1	37
5			5-Chloro-3-ethyl-2-propyl-1H-pyr...	222	C12H15ClN2	1000486-81-4	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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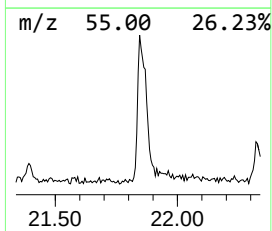
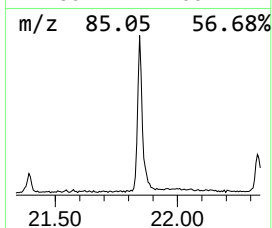
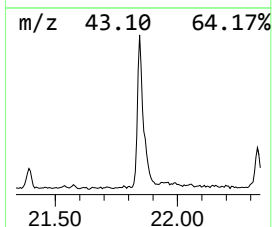
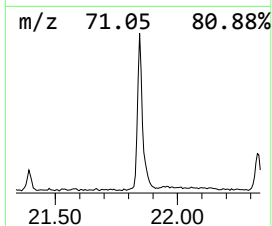
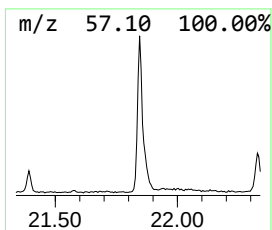
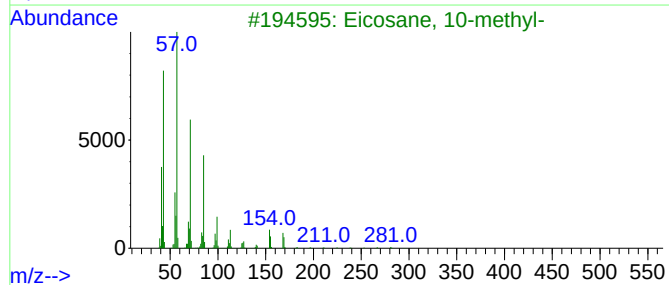
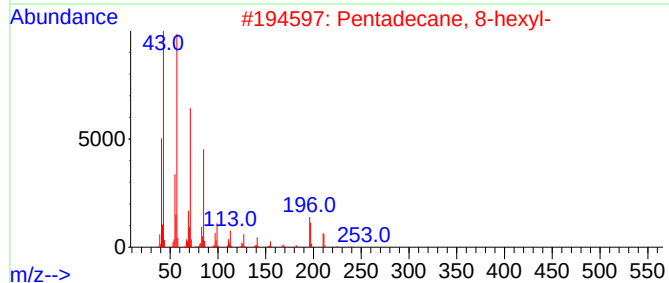
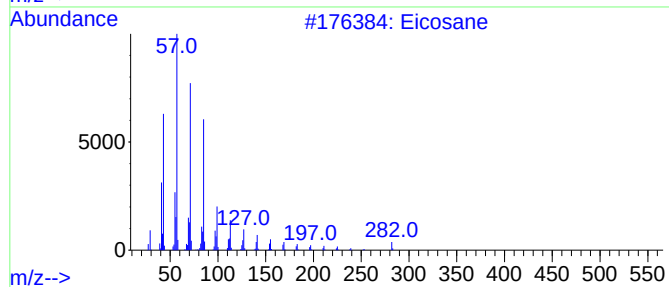
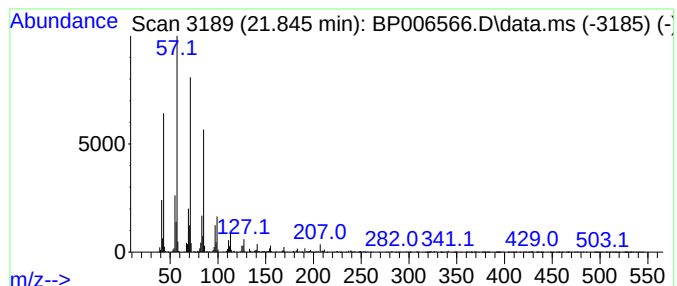
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.840	5.80 ng/ul	699144	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	94
2			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4			Hexacosane	366	C26H54	000630-01-3	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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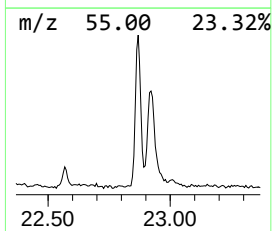
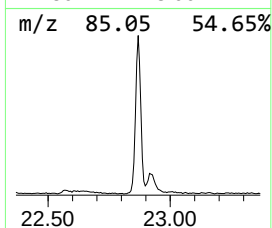
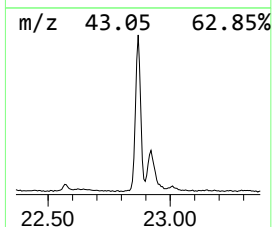
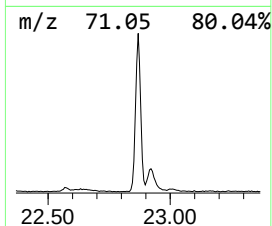
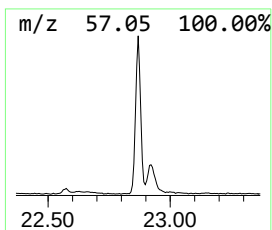
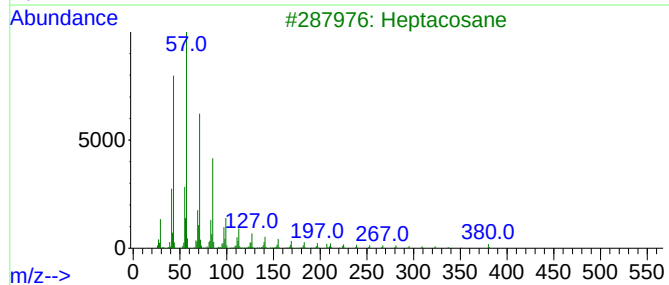
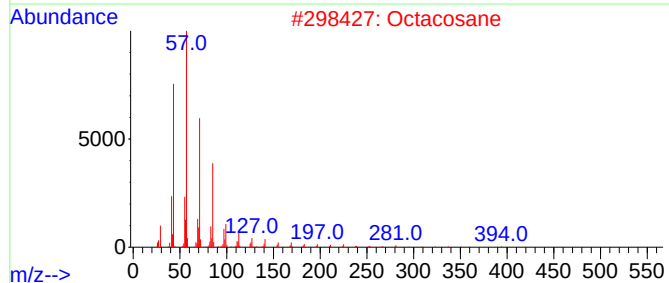
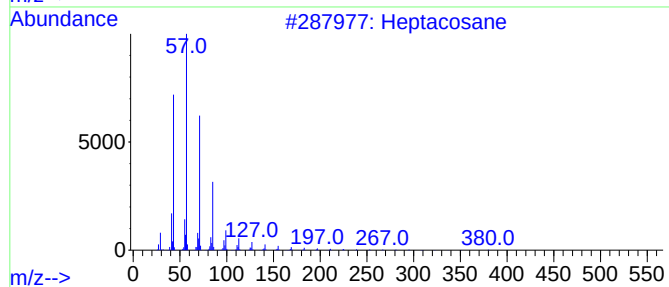
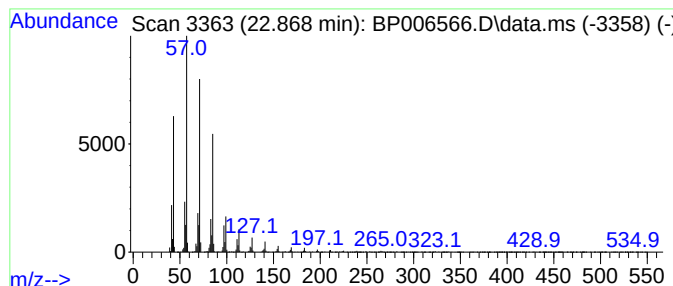
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.870	7.97 ng/ul	1045260	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	94
2			Octacosane	394	C28H58	000630-02-4	94
3			Heptacosane	380	C27H56	000593-49-7	94
4			Heneicosane	296	C21H44	000629-94-7	91
5			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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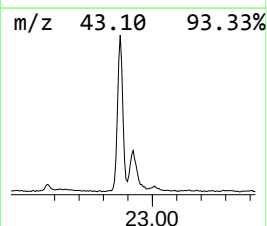
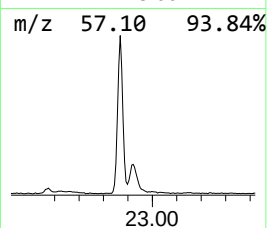
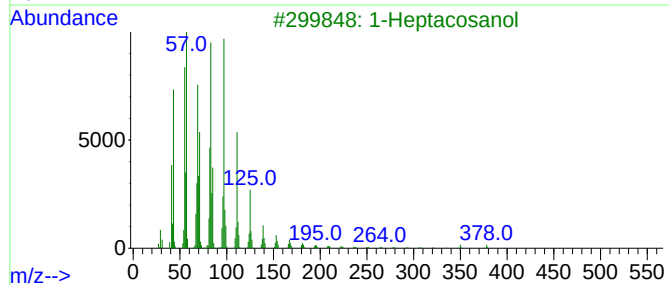
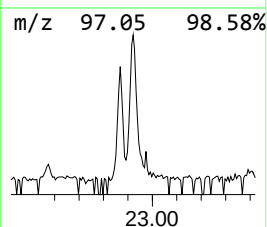
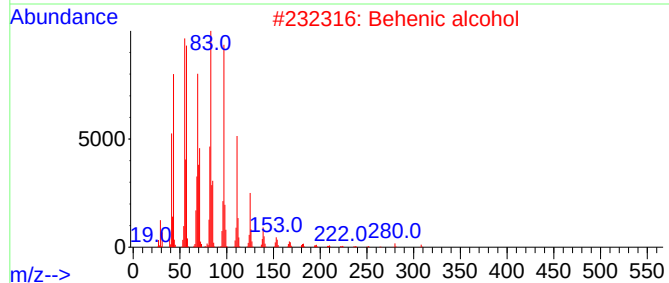
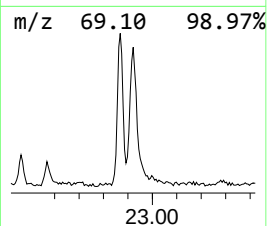
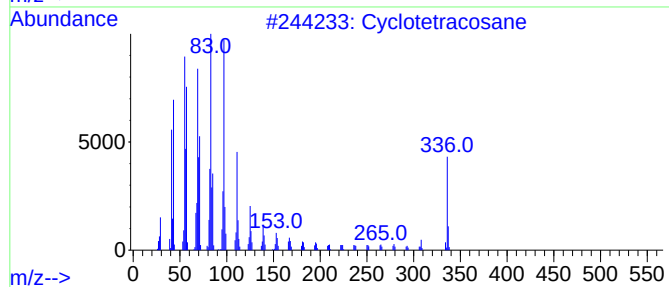
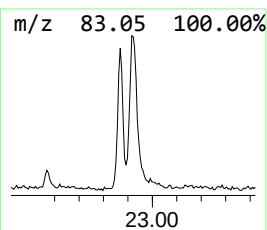
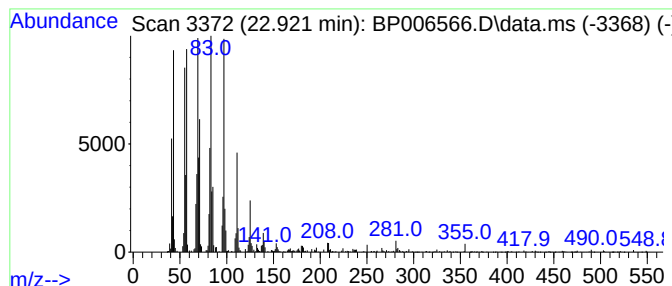
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic22.92 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.920	4.19 ng/ul	548917	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			1-Heptacosanol	396	C27H56O	002004-39-9	93
4			Nonacos-1-ene	406	C29H58	018835-35-3	93
5			1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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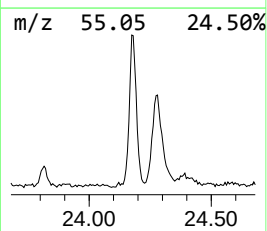
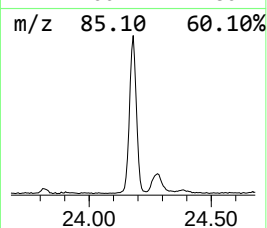
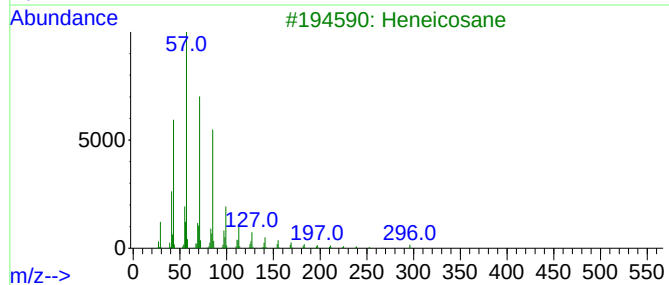
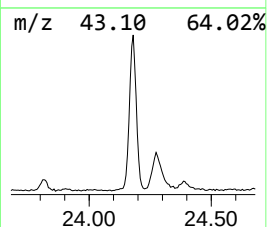
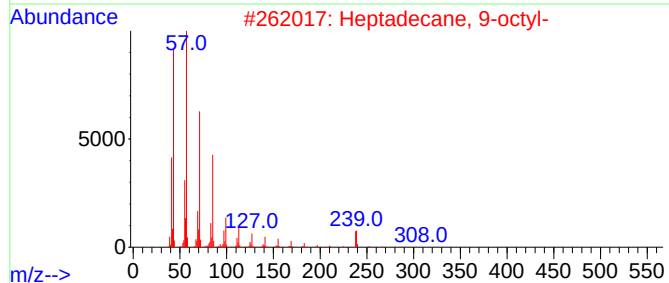
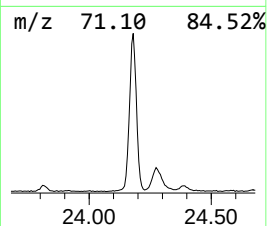
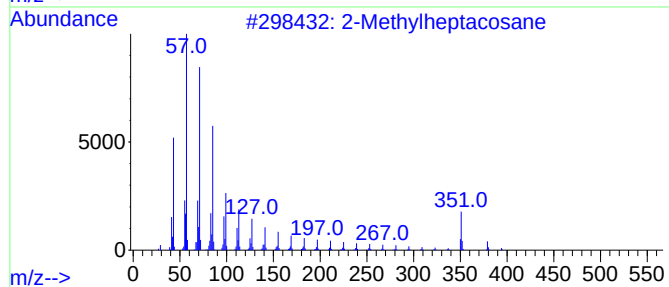
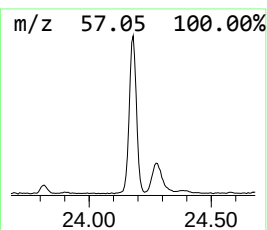
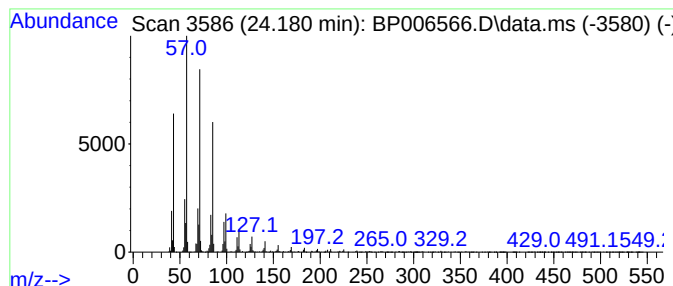
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.180	8.68 ng/ul	1138550	Perylene-d12	23.557

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methylheptacosane	394	C28H58	001561-00-8	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		10-Methylnonadecane	282	C20H42	056862-62-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
Data File : BP006566.D
Acq On : 07 Aug 2021 13:50
Operator : CG/JU
Sample : M3163-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
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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n-Hexadecanoic ...	18.040	2.3	ng/ul	249623	4	17.133	2173730	20.0
2-Dimethylamino...	20.850	3.6	ng/ul	432340	5	21.227	2409310	20.0
(1R,2S,8R,8Ar)-...	20.980	12.9	ng/ul	1548160	5	21.227	2409310	20.0
unknown-01	21.150	5.8	ng/ul	696824	5	21.227	2409310	20.0
(DEL) Alkane: S...	21.840	5.8	ng/ul	699144	5	21.227	2409310	20.0
(DEL) Alkane: S...	22.870	8.0	ng/ul	1045260	6	23.557	2621940	20.0
(DEL) Alkane: C...	22.920	4.2	ng/ul	548917	6	23.557	2621940	20.0
(DEL) Alkane: S...	24.180	8.7	ng/ul	1138550	6	23.557	2621940	20.0