

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
 Data File : BP006614.D
 Acq On : 09 Aug 2021 15:35
 Operator : CG/JU
 Sample : M3208-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C00E1

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-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006614.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.687	98	102	113	rBV	197222	327245	10.71%	0.817%
2	3.781	115	118	124	rVB2	99760	128366	4.20%	0.320%
3	6.934	649	654	668	rVB	611112	991923	32.47%	2.476%
4	7.099	676	682	689	rBV	495145	769660	25.19%	1.921%
5	7.287	708	714	725	rVB	662173	1038126	33.98%	2.592%
6	7.752	786	793	800	rVB	667391	1048873	34.33%	2.618%
7	8.463	908	914	923	rBV	540247	860550	28.17%	2.148%
8	8.575	929	933	938	rVB3	32646	52321	1.71%	0.131%
9	8.911	983	990	997	rBV	607499	1008014	32.99%	2.516%
10	8.975	998	1001	1006	rVB2	38087	51840	1.70%	0.129%
11	9.628	1105	1112	1123	rBV	499661	854194	27.96%	2.132%
12	9.799	1137	1141	1142	rBV4	8222	9849	0.32%	0.025%
13	9.875	1151	1154	1162	rVB8	12905	24747	0.81%	0.062%
14	9.969	1168	1170	1177	rVB8	11871	18355	0.60%	0.046%
15	10.163	1196	1203	1213	rVB	709866	1157144	37.87%	2.889%
16	10.328	1226	1231	1240	rVB	130572	213964	7.00%	0.534%
17	10.534	1257	1266	1272	rBV	905546	1586066	51.91%	3.959%
18	10.587	1272	1275	1282	rVB2	29570	48958	1.60%	0.122%
19	10.681	1282	1291	1303	rBV	488621	917411	30.03%	2.290%
20	11.240	1375	1386	1394	rVB	9622	31089	1.02%	0.078%
21	11.346	1399	1404	1408	rBV6	9091	17531	0.57%	0.044%
22	11.457	1419	1423	1429	rVB4	12409	21312	0.70%	0.053%
23	11.569	1437	1442	1447	rBV	45725	72125	2.36%	0.180%
24	11.816	1478	1484	1487	rVB8	5450	10768	0.35%	0.027%
25	11.969	1503	1510	1516	rBV	58825	89481	2.93%	0.223%
26	12.046	1519	1523	1525	rVV5	6587	8811	0.29%	0.022%
27	12.122	1528	1536	1541	rVV4	15392	39972	1.31%	0.100%
28	12.199	1544	1549	1557	rVB	54356	86160	2.82%	0.215%
29	12.334	1568	1572	1578	rBV8	8093	17851	0.58%	0.045%
30	12.416	1582	1586	1591	rVV	31649	51555	1.69%	0.129%
31	12.493	1595	1599	1603	rBV7	8883	13876	0.45%	0.035%
32	12.604	1614	1618	1622	rBV7	4242	7837	0.26%	0.020%
33	12.663	1622	1628	1633	rVV7	16930	37917	1.24%	0.095%
34	12.740	1635	1641	1644	rVV8	10309	27264	0.89%	0.068%
35	12.787	1645	1649	1657	rVV5	13728	33167	1.09%	0.083%
36	12.869	1659	1663	1665	rVV5	9722	13815	0.45%	0.034%

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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 >

Peak separation: 5

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

37	12.928	1668	1673	1677	rVV2	26469	46238	1.51%	0.115%
38	12.975	1679	1681	1688	rVB8	8029	16130	0.53%	0.040%
39	13.216	1716	1722	1729	rVB2	103475	147381	4.82%	0.368%
40	13.316	1737	1739	1744	rVB6	5439	8639	0.28%	0.022%

41	13.416	1753	1756	1759	rBV5	5287	9221	0.30%	0.023%
42	13.546	1771	1778	1779	rBV	22216	28716	0.94%	0.072%
43	13.563	1779	1781	1786	rVB3	21700	27408	0.90%	0.068%
44	13.663	1794	1798	1800	rBV5	13406	17080	0.56%	0.043%
45	13.704	1800	1805	1809	rBV3	37325	57620	1.89%	0.144%

46	13.757	1809	1814	1817	rBV2	29460	44866	1.47%	0.112%
47	13.804	1817	1822	1828	rVB	1256126	1679299	54.96%	4.192%
48	13.875	1830	1834	1838	rVB2	33064	44989	1.47%	0.112%
49	13.916	1838	1841	1842	rBV3	14410	15837	0.52%	0.040%
50	14.075	1858	1868	1880	rBV	1500854	2197602	71.93%	5.486%

51	14.210	1888	1891	1896	rVB7	11539	15083	0.49%	0.038%
52	14.293	1900	1905	1911	rVB	109740	137926	4.51%	0.344%
53	14.381	1911	1920	1928	rBV	1225749	1906343	62.39%	4.759%
54	14.446	1928	1931	1933	rBV4	7080	9766	0.32%	0.024%
55	14.475	1933	1936	1940	rVB3	18408	25910	0.85%	0.065%

56	14.604	1951	1958	1969	rBV	431276	759457	24.86%	1.896%
57	14.751	1977	1983	1986	rBV4	35000	63924	2.09%	0.160%
58	14.787	1986	1989	1992	rVB3	28037	38875	1.27%	0.097%
59	14.863	1998	2002	2006	rVB3	28128	42024	1.38%	0.105%
60	14.916	2006	2011	2018	rVB5	18181	34126	1.12%	0.085%

61	15.004	2019	2026	2031	rBV4	29072	57112	1.87%	0.143%
62	15.057	2031	2035	2040	rVB	34831	50026	1.64%	0.125%
63	15.140	2046	2049	2052	rBV	17765	23889	0.78%	0.060%
64	15.175	2052	2055	2059	rVV4	17803	22789	0.75%	0.057%
65	15.245	2063	2067	2072	rVB	117197	144556	4.73%	0.361%

66	15.316	2076	2079	2082	rBV4	10726	12252	0.40%	0.031%
67	15.381	2083	2090	2096	rBV	1764975	2471672	80.90%	6.170%
68	15.440	2097	2100	2104	rVV5	27964	46459	1.52%	0.116%
69	15.510	2107	2112	2117	rVB	232149	324949	10.64%	0.811%
70	15.622	2129	2131	2133	rVV3	10788	8882	0.29%	0.022%

71	15.651	2133	2136	2141	rVB3	27679	42452	1.39%	0.106%
72	15.728	2146	2149	2157	rVB4	29069	51548	1.69%	0.129%
73	15.810	2160	2163	2166	rBV4	13800	18121	0.59%	0.045%
74	15.963	2183	2189	2195	rBV5	25757	62013	2.03%	0.155%
75	16.116	2210	2215	2222	rBV	159117	266202	8.71%	0.665%

76	16.204	2228	2230	2235	rBV6	12325	21181	0.69%	0.053%
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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 >
 Peak separation: 5

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

77	16.492	2276	2279	2285	rVB2	18580	27553	0.90%	0.069%
78	16.687	2307	2312	2314	rBV4	30659	54281	1.78%	0.136%
79	16.716	2315	2317	2320	rVB3	25663	25141	0.82%	0.063%
80	16.792	2326	2330	2338	rBV3	56513	101125	3.31%	0.252%
81	16.869	2339	2343	2346	rBV2	38568	62881	2.06%	0.157%
82	16.910	2346	2350	2355	rVB	132483	169987	5.56%	0.424%
83	17.140	2383	2389	2393	rBV	1429105	2018557	66.07%	5.039%
84	17.181	2393	2396	2400	rVV	252841	335557	10.98%	0.838%
85	17.240	2400	2406	2416	rVV	1698997	2548167	83.40%	6.361%
86	17.457	2440	2443	2447	rVB	53214	71267	2.33%	0.178%
87	17.651	2472	2476	2480	rBV	135054	165772	5.43%	0.414%
88	17.828	2503	2506	2510	rVB2	37431	40037	1.31%	0.100%
89	18.010	2534	2537	2539	rBV	77430	88118	2.88%	0.220%
90	18.051	2539	2544	2552	rVB2	214252	417322	13.66%	1.042%
91	18.192	2565	2568	2572	rVB2	108463	126734	4.15%	0.316%
92	18.322	2587	2590	2598	rBV2	150417	289327	9.47%	0.722%
93	18.498	2617	2620	2625	rBV	112367	152251	4.98%	0.380%
94	18.939	2692	2695	2701	rBV2	134033	226497	7.41%	0.565%
95	19.157	2728	2732	2736	rBV	570427	742393	24.30%	1.853%
96	19.481	2782	2787	2791	rBV	2172019	3055318	100.00%	7.627%
97	20.963	3036	3039	3044	rBV3	437925	505819	16.56%	1.263%
98	21.239	3082	3086	3089	rVB	1513334	1966318	64.36%	4.909%
99	23.422	3453	3457	3463	rVB2	1229776	2077619	68.00%	5.187%
100	23.574	3478	3483	3490	rVB2	1106248	2103131	68.84%	5.250%

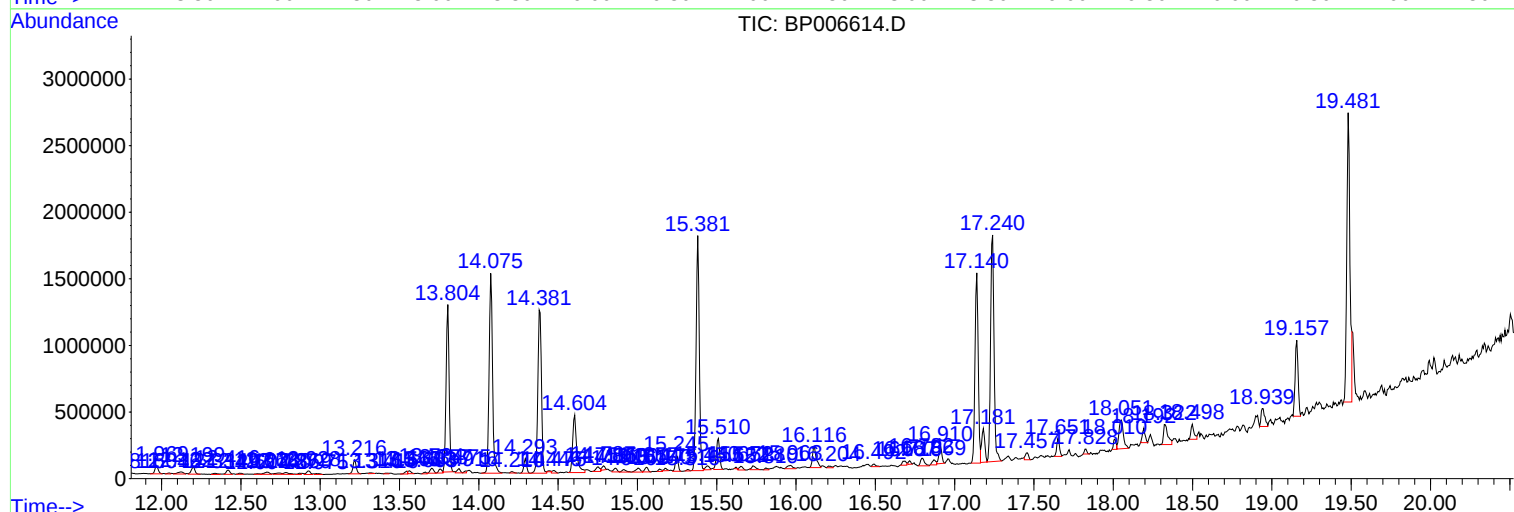
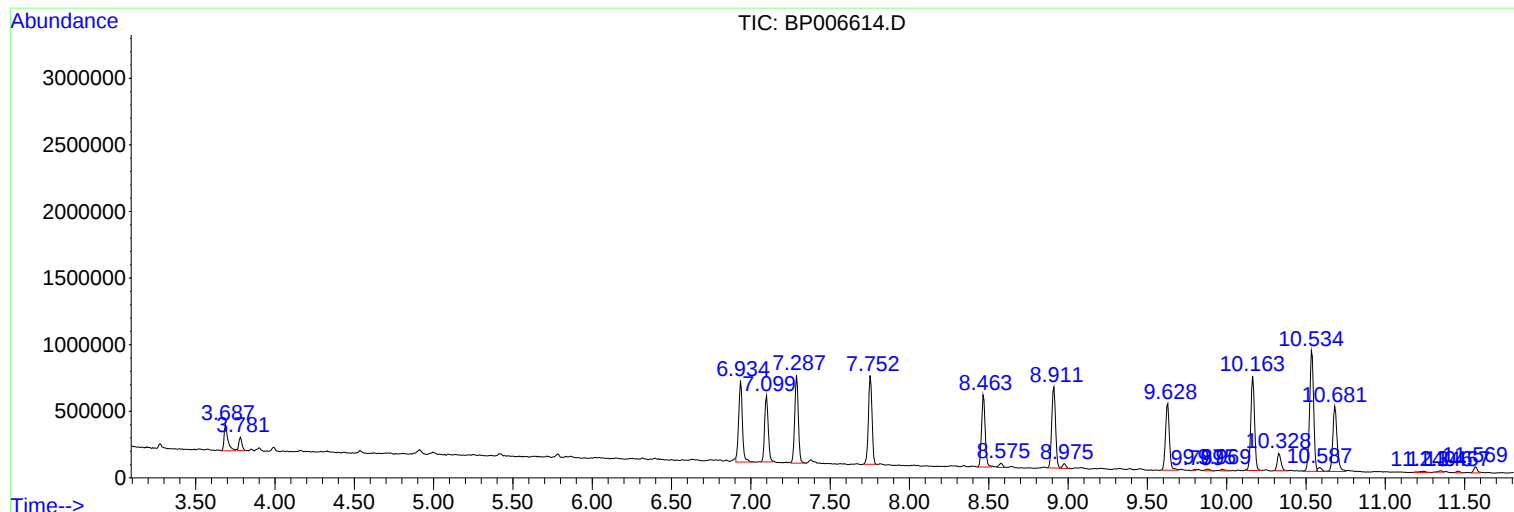
Sum of corrected areas: 40057872

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Instrument :
BNA_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



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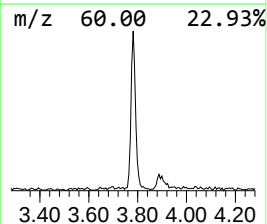
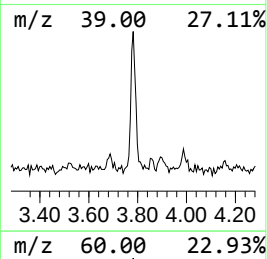
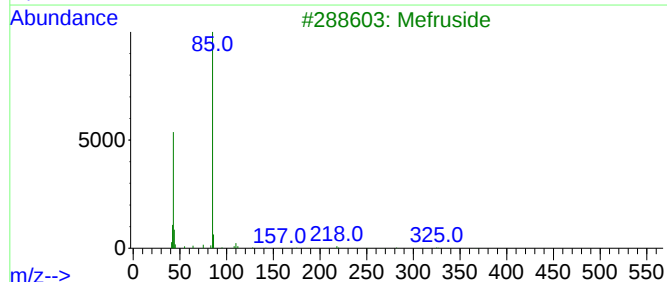
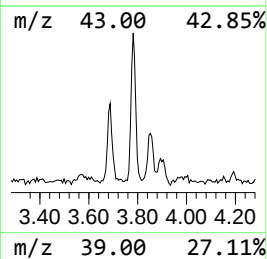
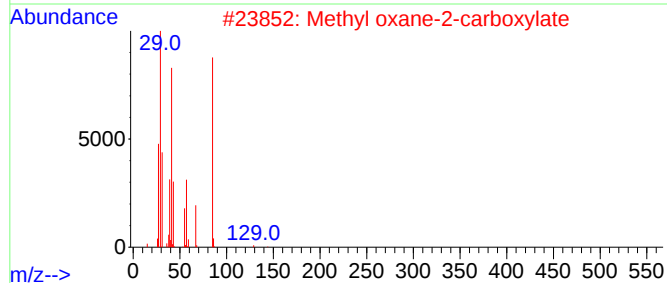
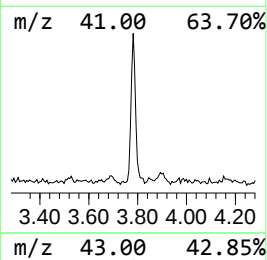
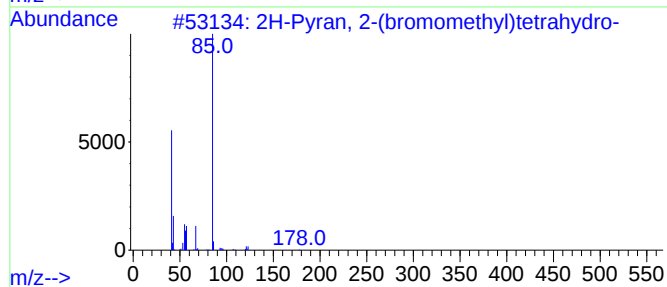
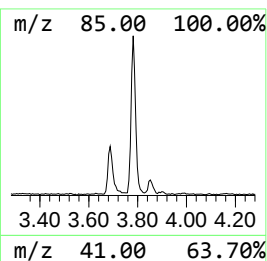
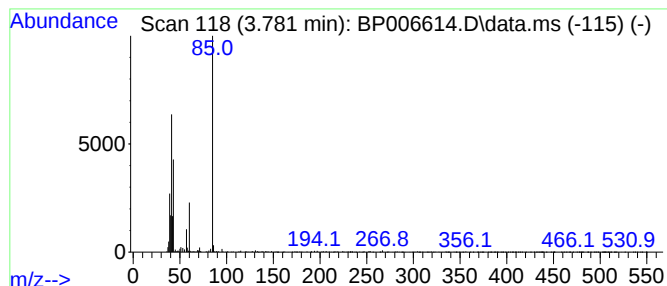
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.780	2.45 ng/ul	128366	1,4-Dichlorobenzene-d4	7.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, 2-(bromomethyl)tetrahy...	178	C6H11BrO	034723-82-5	40		
2	Methyl oxane-2-carboxylate	144	C7H12O3	105441-86-9	38		
3	Mefruside	382	C13H19ClN2O5S2	007195-27-9	33		
4	2(3H)-Furanone, 5-butylidihydro-	142	C8H14O2	000104-50-7	9		
5	5-(Bromomethyl)oxolan-2-one	178	C5H7BrO2	099393-06-3	9		



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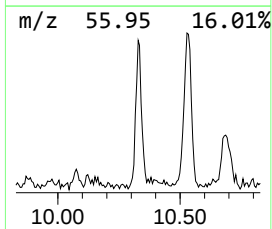
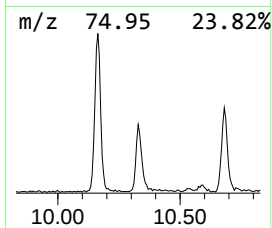
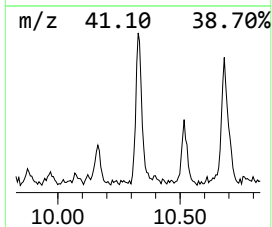
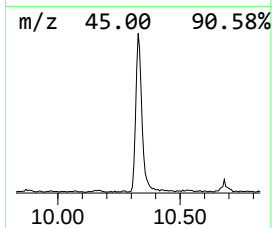
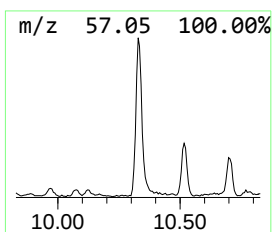
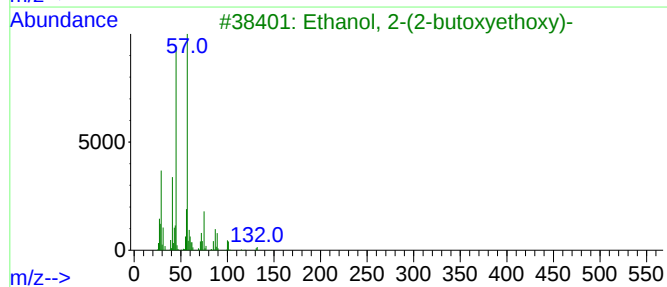
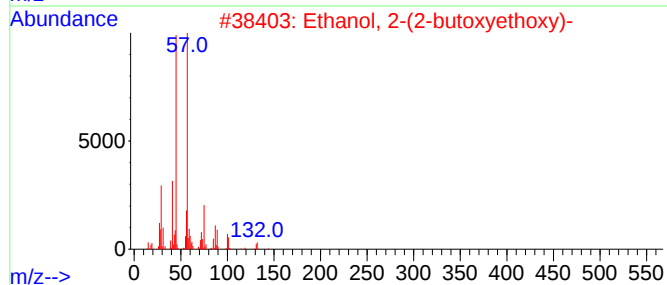
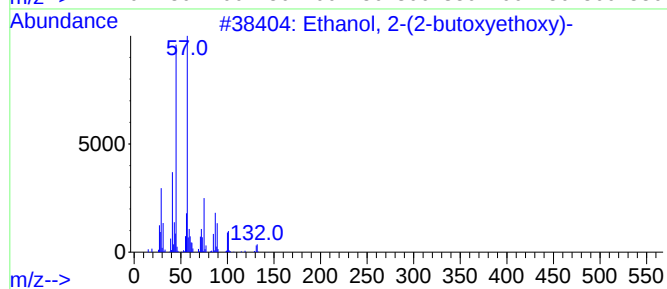
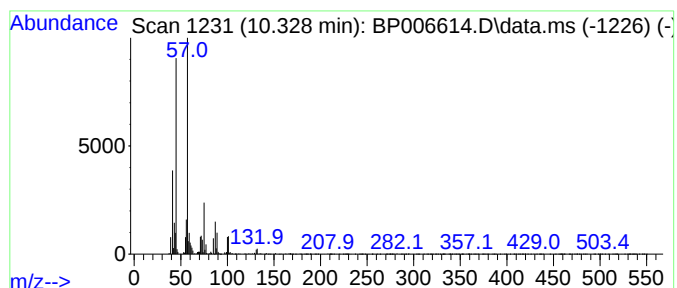
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.330	2.70 ng/ul	213964	Naphthalene-d8	10.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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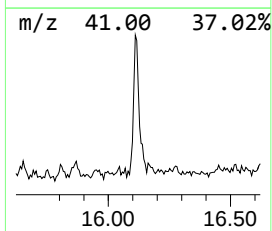
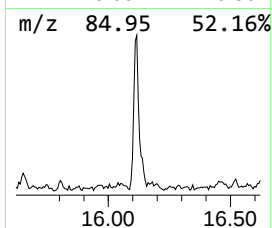
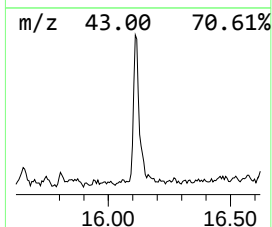
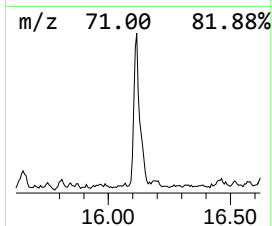
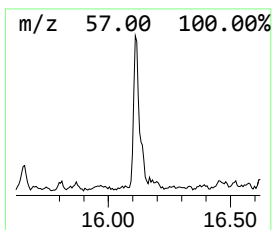
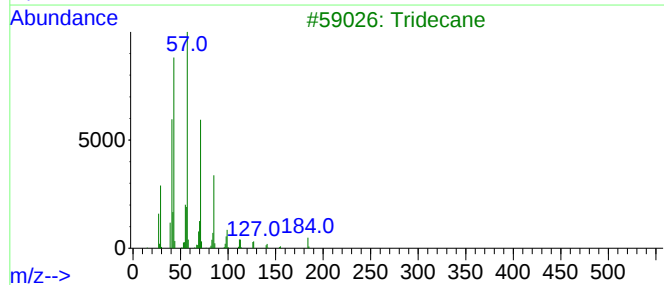
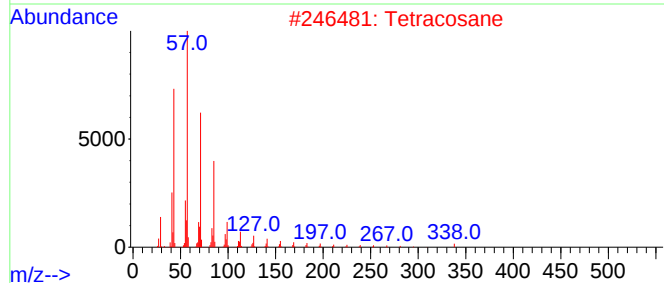
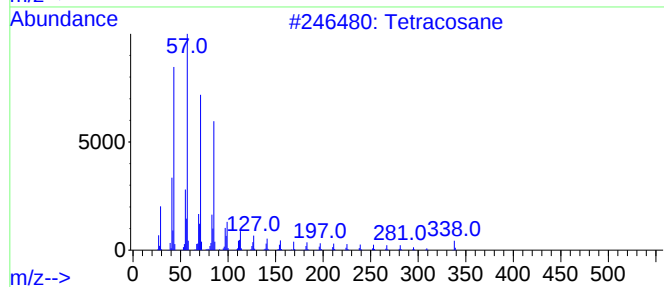
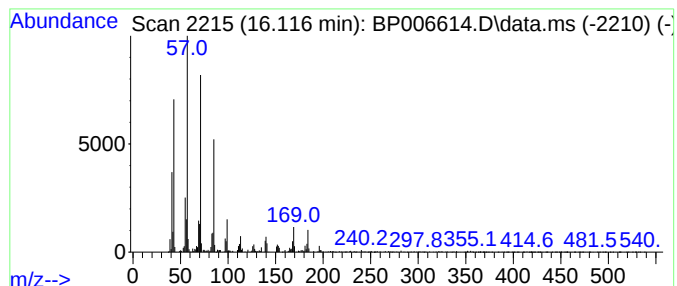
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.120	2.64 ng/ul	266202	Phenanthrene-d10	17.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	83
2			Tetracosane	338	C24H50	000646-31-1	83
3			Tridecane	184	C13H28	000629-50-5	83
4			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006614.D
Acq On : 09 Aug 2021 15:35
Operator : CG/JU
Sample : M3208-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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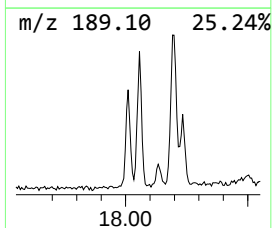
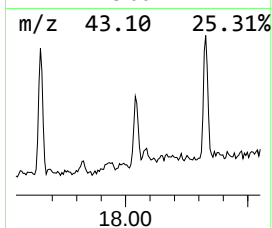
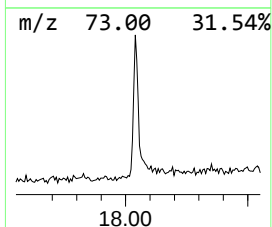
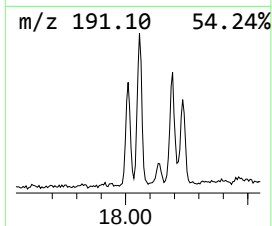
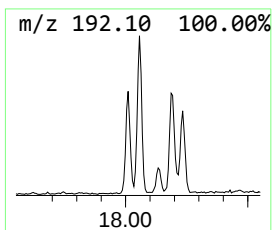
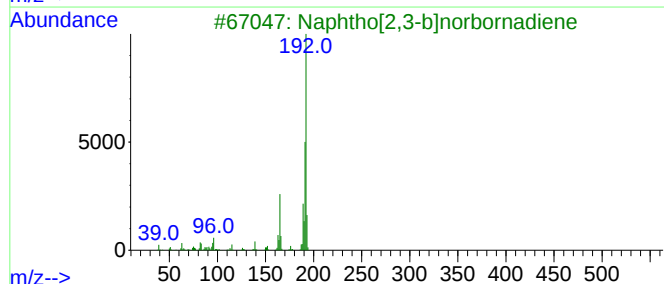
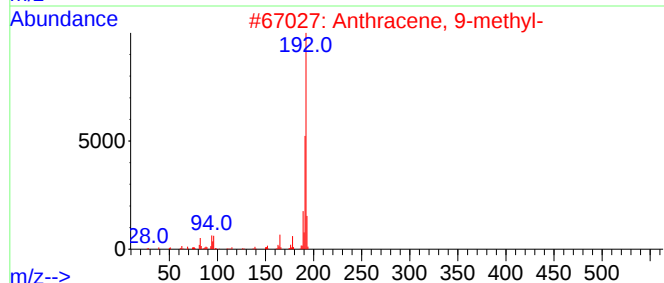
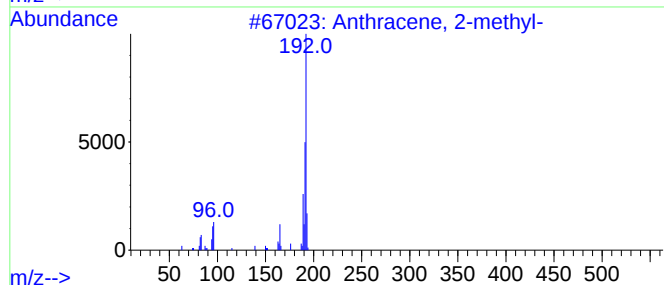
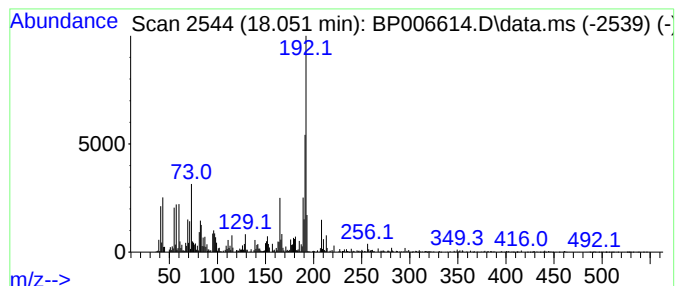
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.050	4.13 ng/ul	417322	Phenanthrene-d10	17.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	86
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
3			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	83
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006614.D
Acq On : 09 Aug 2021 15:35
Operator : CG/JU
Sample : M3208-01
Misc :
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Instrument :
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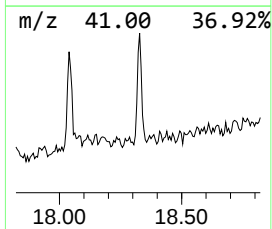
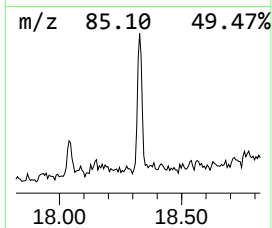
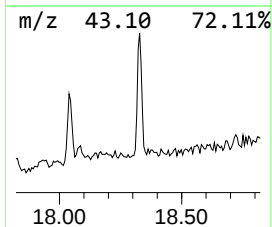
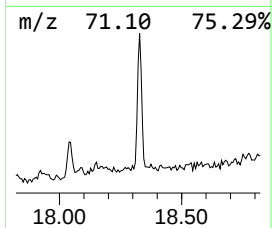
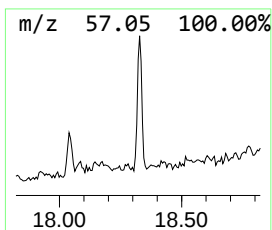
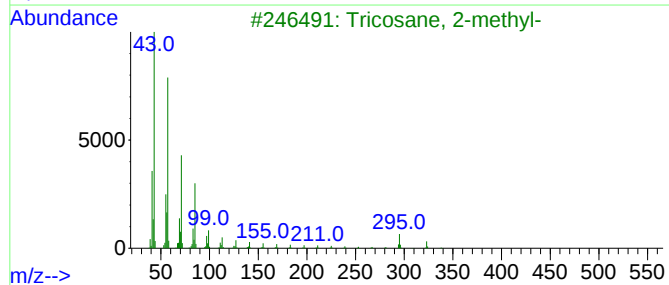
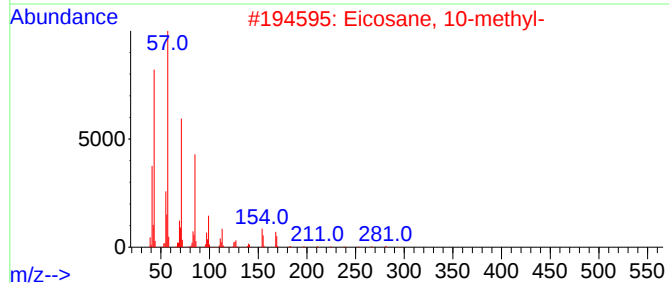
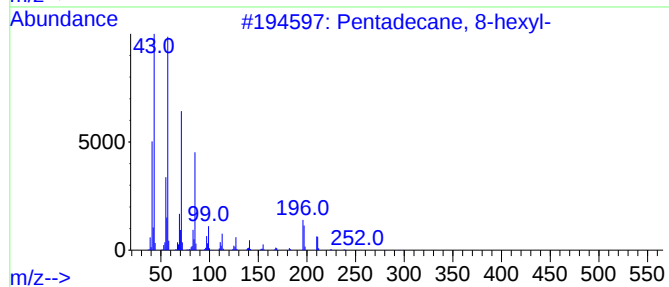
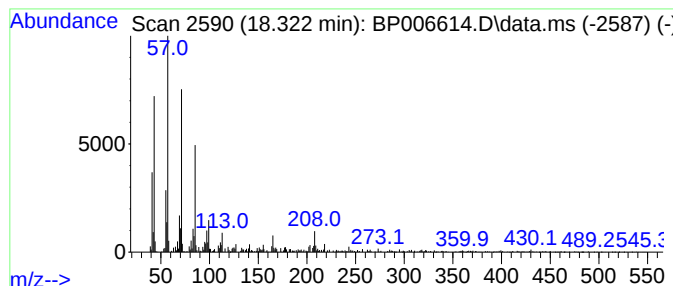
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.320	2.87 ng/ul	289327	Phenanthrene-d10	17.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	87
3			Tricosane, 2-methyl-	338	C24H50	001928-30-9	87
4			20.Xi.-Lanosta-7,9(11)-diene-3.b...	458	C30H50O3	025116-58-9	84
5			Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006614.D
Acq On : 09 Aug 2021 15:35
Operator : CG/JU
Sample : M3208-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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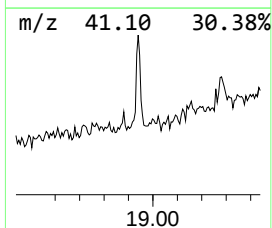
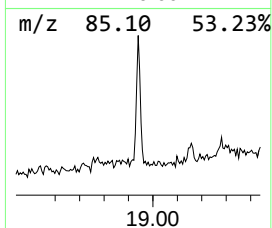
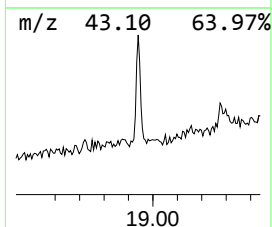
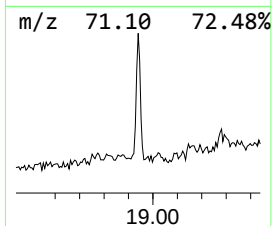
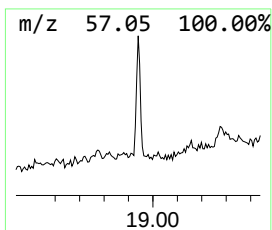
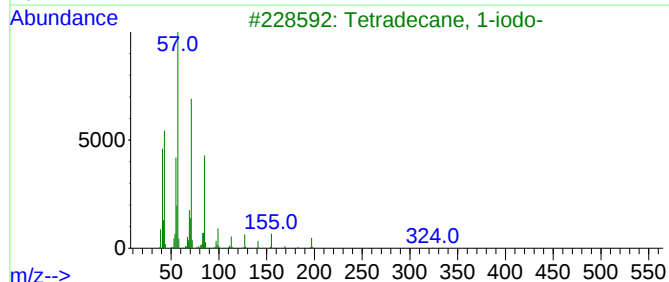
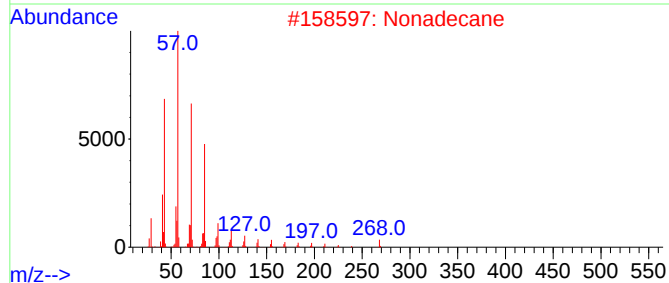
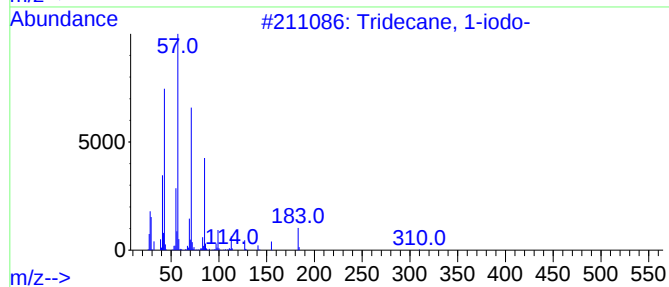
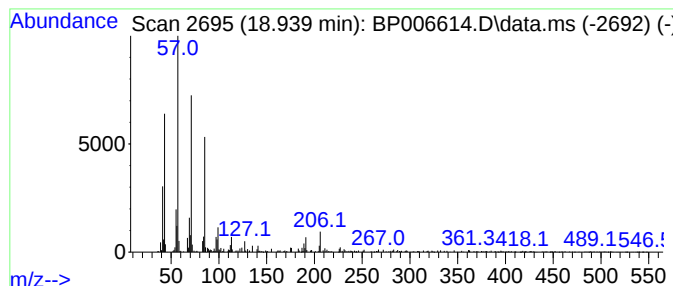
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tridecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.940	2.24 ng/ul	226497	Phenanthrene-d10	17.140

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
2			Nonadecane	268	C19H40	000629-92-5	74
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
4			Tetradecane	198	C14H30	000629-59-4	72
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006614.D
Acq On : 09 Aug 2021 15:35
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Sample : M3208-01
Misc :
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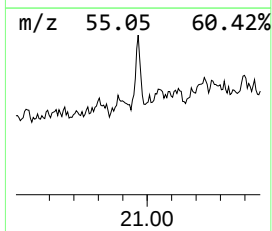
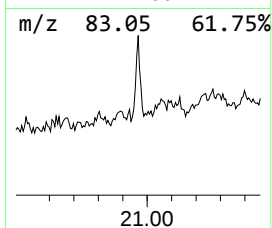
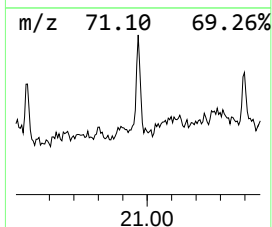
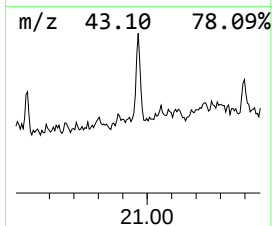
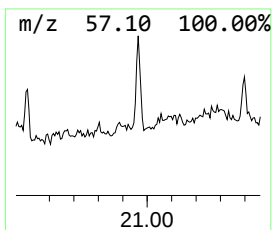
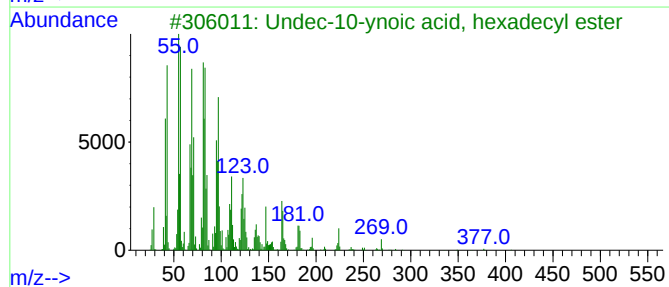
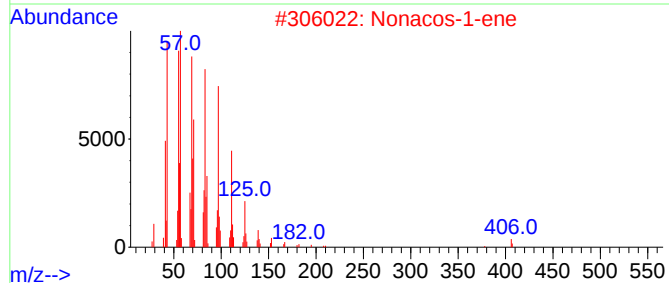
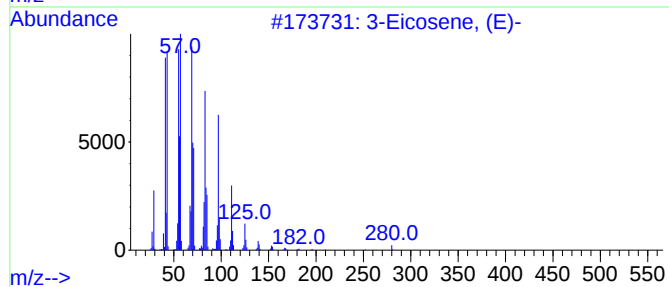
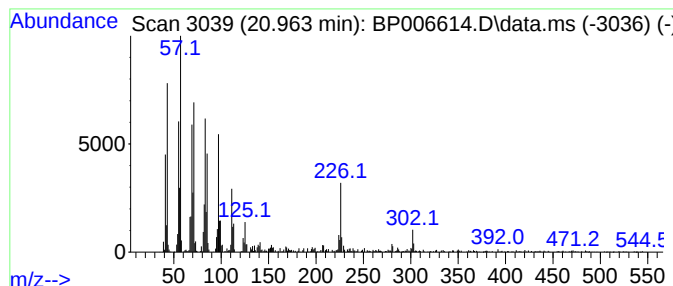
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.960	5.14 ng/ul	505819	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-			280	C20H40	074685-33-9	95
2	Nonacos-1-ene			406	C29H58	018835-35-3	83
3	Undec-10-ynoic acid, hexadecyl e...			406	C27H50O2	1000406-16-9	83
4	Cyclohexadecane, 1,2-diethyl-			280	C20H40	1000155-85-3	78
5	Ethanol, 2-(tetradecyloxy)-			258	C16H34O2	002136-70-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006614.D
Acq On : 09 Aug 2021 15:35
Operator : CG/JU
Sample : M3208-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00E1

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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Ethanol, 2-(2-b...	10.330	2.7	ng/ul	213964	2	10.534	1586070	20.0
(DEL) Alkane: S...	16.120	2.6	ng/ul	266202	4	17.140	2018560	20.0
Anthracene, 2-m...	18.050	4.1	ng/ul	417322	4	17.140	2018560	20.0
(DEL) Alkane: S...	18.320	2.9	ng/ul	289327	4	17.140	2018560	20.0
Tridecane, 1-iodo-	18.940	2.2	ng/ul	226497	4	17.140	2018560	20.0
(DEL) Alkane: S...	20.960	5.1	ng/ul	505819	5	21.239	1966320	20.0