

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007631.D
 Acq On : 23 Oct 2021 15:33
 Operator : CG/JU
 Sample : M4262-16
 Misc :
 ALS Vial : 91 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY49

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007631.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.376	11	15	18	rBV	43273	59147	1.10%	0.093%
2	3.411	18	21	27	rVB	61190	79404	1.47%	0.125%
3	3.658	59	63	69	rBV3	16017	28016	0.52%	0.044%
4	3.723	69	74	79	rVB	22050	33193	0.61%	0.052%
5	3.793	82	86	100	rBV	352276	562747	10.42%	0.884%
6	4.034	122	127	132	rVB	25911	40518	0.75%	0.064%
7	4.123	139	142	147	rVB	17465	21813	0.40%	0.034%
8	4.217	154	158	164	rBV5	11474	16480	0.31%	0.026%
9	5.052	293	300	306	rBV6	11545	20024	0.37%	0.031%
10	5.787	418	425	429	rBV	14489	25404	0.47%	0.040%
11	5.828	429	432	438	rVB2	16349	25494	0.47%	0.040%
12	6.022	462	465	471	rVB	11108	16567	0.31%	0.026%
13	6.328	511	517	520	rBV3	8332	14627	0.27%	0.023%
14	6.422	527	533	540	rBV	17348	30839	0.57%	0.048%
15	6.952	614	623	628	rBV8	12068	34049	0.63%	0.053%
16	7.111	643	650	659	rVB	250053	410510	7.60%	0.645%
17	7.269	671	677	692	rVB	624558	1009745	18.69%	1.586%
18	7.475	703	712	720	rBV	661813	1038501	19.23%	1.631%
19	7.546	720	724	728	rBV	15846	22040	0.41%	0.035%
20	7.946	784	792	799	rBV	817288	1326837	24.57%	2.084%
21	8.011	799	803	812	rVB2	17480	31609	0.59%	0.050%
22	8.187	824	833	842	rBV3	17828	39322	0.73%	0.062%
23	8.469	874	881	883	rBV	9322	15567	0.29%	0.024%
24	8.505	883	887	891	rVB2	10931	18975	0.35%	0.030%
25	8.652	905	912	917	rBV	545578	879940	16.29%	1.382%
26	8.711	917	922	929	rVV	653431	1003527	18.58%	1.576%
27	9.099	980	988	998	rBV	758465	1243022	23.01%	1.953%
28	9.275	1014	1018	1023	rVV4	14218	18486	0.34%	0.029%
29	9.563	1057	1067	1075	rVB2	38802	81301	1.51%	0.128%
30	9.822	1102	1111	1124	rBV	567191	945596	17.51%	1.485%
31	10.069	1148	1153	1162	rVB5	14439	34474	0.64%	0.054%
32	10.169	1164	1170	1173	rBV7	8741	18770	0.35%	0.029%
33	10.369	1195	1204	1213	rBV	986739	1621740	30.03%	2.548%
34	10.528	1223	1231	1247	rBV	949955	1503967	27.85%	2.363%
35	10.746	1259	1268	1276	rBV	1092254	1848011	34.21%	2.903%
36	10.881	1283	1291	1305	rBV	907004	1569271	29.05%	2.465%

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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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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
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37	11.099	1322	1328	1333	rBV9	7252	16480	0.31%	0.026%
38	11.740	1432	1437	1441	rVV8	7414	14919	0.28%	0.023%
39	11.946	1466	1472	1475	rBV2	17564	28335	0.52%	0.045%
40	11.987	1475	1479	1482	rVV2	19138	31507	0.58%	0.049%
41	12.057	1486	1491	1497	rVB4	24035	41173	0.76%	0.065%
42	12.222	1515	1519	1531	rVV3	37910	95709	1.77%	0.150%
43	12.328	1531	1537	1543	rVV4	32729	97046	1.80%	0.152%
44	12.387	1543	1547	1553	rVV7	26891	61327	1.14%	0.096%
45	12.434	1553	1555	1562	rVB8	12482	22520	0.42%	0.035%
46	12.828	1617	1622	1625	rBV6	18438	34560	0.64%	0.054%
47	12.952	1639	1643	1648	rVB3	13850	21649	0.40%	0.034%
48	13.199	1681	1685	1691	rVB2	26035	36519	0.68%	0.057%
49	13.363	1708	1713	1716	rBV2	20495	32935	0.61%	0.052%
50	13.428	1720	1724	1730	rVV6	17622	29337	0.54%	0.046%
51	13.487	1730	1734	1743	rVV6	17828	37101	0.69%	0.058%
52	13.710	1766	1772	1778	rBV2	15009	29035	0.54%	0.046%
53	13.987	1813	1819	1825	rVV	1807789	2566820	47.52%	4.032%
54	14.040	1826	1828	1831	rVV4	15058	17282	0.32%	0.027%
55	14.269	1854	1867	1876	rBV	1962601	3004177	55.62%	4.719%
56	14.428	1890	1894	1899	rBV7	31835	53425	0.99%	0.084%
57	14.498	1900	1906	1907	rVV5	29044	45148	0.84%	0.071%
58	14.528	1907	1911	1914	rVV3	80267	138762	2.57%	0.218%
59	14.581	1914	1920	1932	rVB	1763396	2826080	52.32%	4.440%
60	14.775	1948	1953	1961	rVB2	154618	233402	4.32%	0.367%
61	14.940	1969	1981	1988	rBV2	56081	205997	3.81%	0.324%
62	15.010	1988	1993	1995	rBV3	47161	75065	1.39%	0.118%
63	15.157	2015	2018	2019	rBV2	19211	17615	0.33%	0.028%
64	15.328	2039	2047	2054	rVB2	59858	136705	2.53%	0.215%
65	15.393	2054	2058	2062	rBV3	58323	91335	1.69%	0.143%
66	15.440	2062	2066	2071	rVV	215513	283752	5.25%	0.446%
67	15.498	2071	2076	2082	rVV5	33138	63591	1.18%	0.100%
68	15.575	2082	2089	2097	rVB	2695689	3873763	71.72%	6.085%
69	15.687	2102	2108	2116	rBV	256456	361566	6.69%	0.568%
70	15.816	2123	2130	2135	rVB	30013	46238	0.86%	0.073%
71	15.945	2145	2152	2157	rBV9	31368	60637	1.12%	0.095%
72	16.022	2161	2165	2170	rVB5	24042	38843	0.72%	0.061%
73	16.140	2180	2185	2186	rBV5	13964	21964	0.41%	0.035%
74	16.216	2195	2198	2201	rBV4	16712	23107	0.43%	0.036%
75	16.310	2211	2214	2219	rVB3	21397	25684	0.48%	0.040%
76	16.363	2220	2223	2227	rVB3	22887	30108	0.56%	0.047%

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77	16.740	2283	2287	2296	rBV2	101027	147835	2.74%	0.232%
78	16.992	2327	2330	2337	rBV7	10911	19536	0.36%	0.031%
79	17.275	2374	2378	2381	rBV6	11230	21465	0.40%	0.034%
80	17.334	2382	2388	2394	rVV	2191759	3056341	56.59%	4.801%
81	17.434	2399	2405	2412	rVV2	3046628	4465596	82.68%	7.015%
82	17.604	2431	2434	2435	rBV2	15115	15011	0.28%	0.024%
83	17.728	2452	2455	2461	rVB3	47354	63812	1.18%	0.100%
84	17.875	2475	2480	2484	rBV	340516	405457	7.51%	0.637%
85	17.916	2484	2487	2494	rVB	166223	215225	3.98%	0.338%
86	18.016	2499	2504	2511	rVB	191005	228030	4.22%	0.358%
87	18.234	2536	2541	2550	rBV	1892383	2652260	49.10%	4.166%
88	18.310	2550	2554	2558	rVB3	156825	238479	4.42%	0.375%
89	19.063	2678	2682	2688	rBV5	41477	83838	1.55%	0.132%
90	19.145	2693	2696	2699	rVB3	51099	54953	1.02%	0.086%
91	19.245	2709	2713	2716	rBV2	42421	58165	1.08%	0.091%
92	19.316	2721	2725	2727	rBV	68188	99798	1.85%	0.157%
93	19.463	2746	2750	2760	rBV2	285782	388204	7.19%	0.610%
94	19.669	2779	2785	2792	rBV	3856602	5292658	97.99%	8.314%
95	19.834	2811	2813	2819	rVB6	36145	42289	0.78%	0.066%
96	21.133	3030	3034	3041	rBV	819948	1076850	19.94%	1.692%
97	21.422	3077	3083	3088	rBV	2658353	3579468	66.27%	5.623%
98	23.710	3464	3472	3480	rVB2	2669687	5401204	100.00%	8.485%
99	23.869	3492	3499	3507	rVB2	1744766	3690151	68.32%	5.797%
100	27.421	4095	4103	4112	rBV5	555898	1799690	33.32%	2.827%

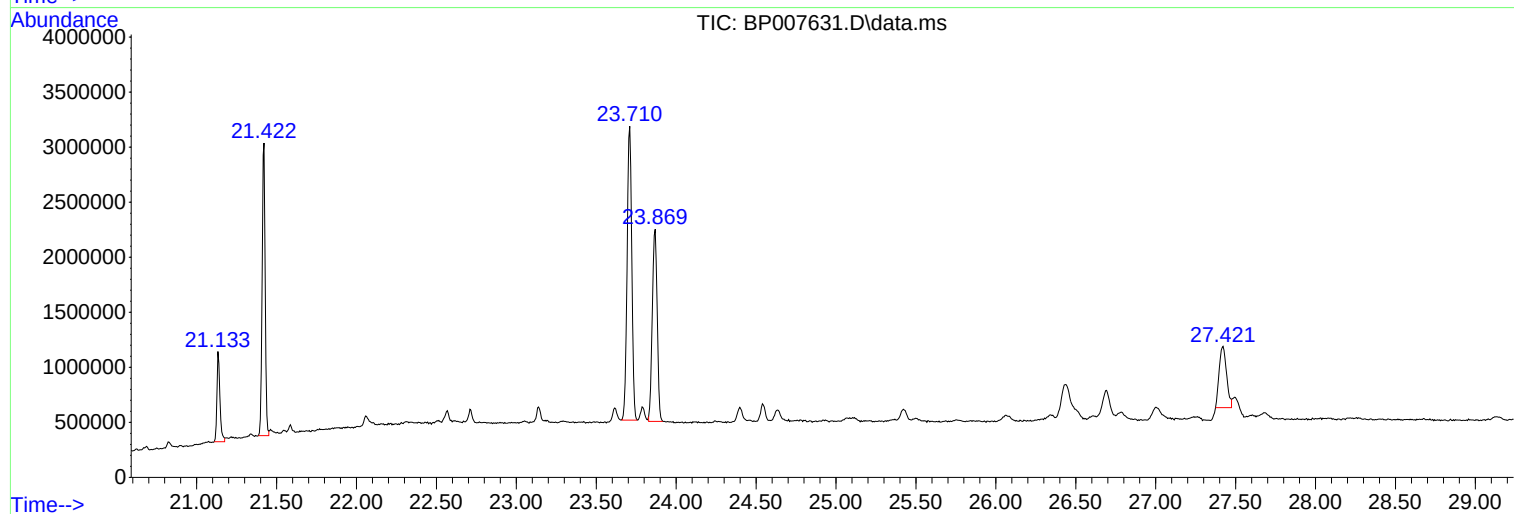
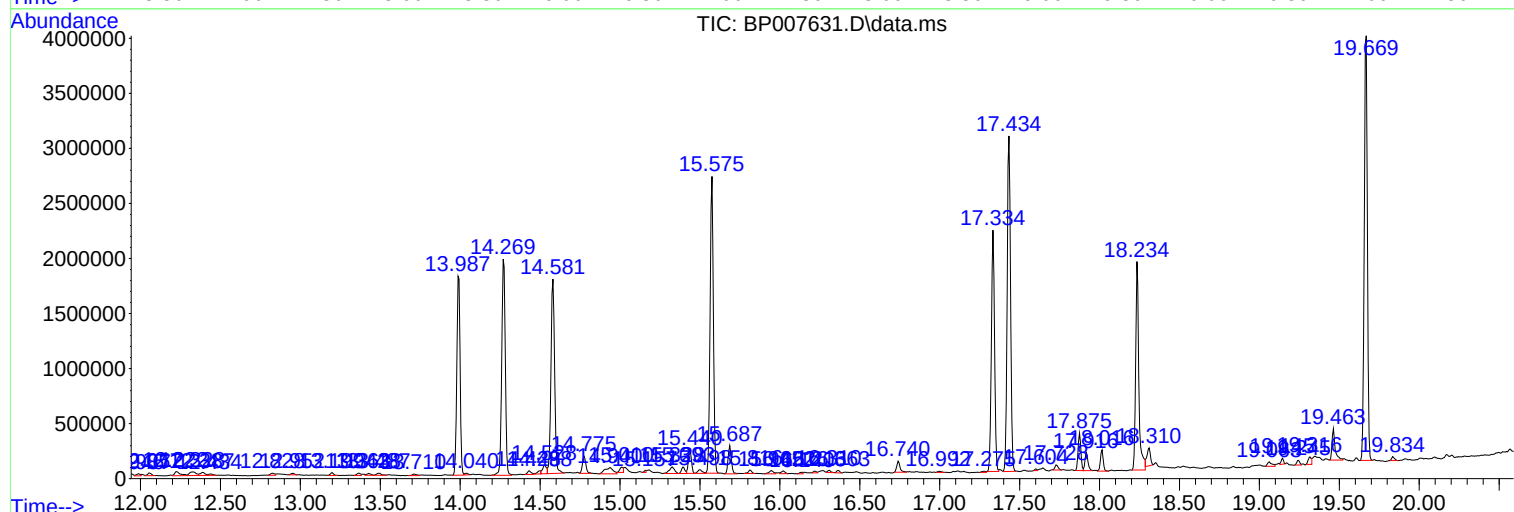
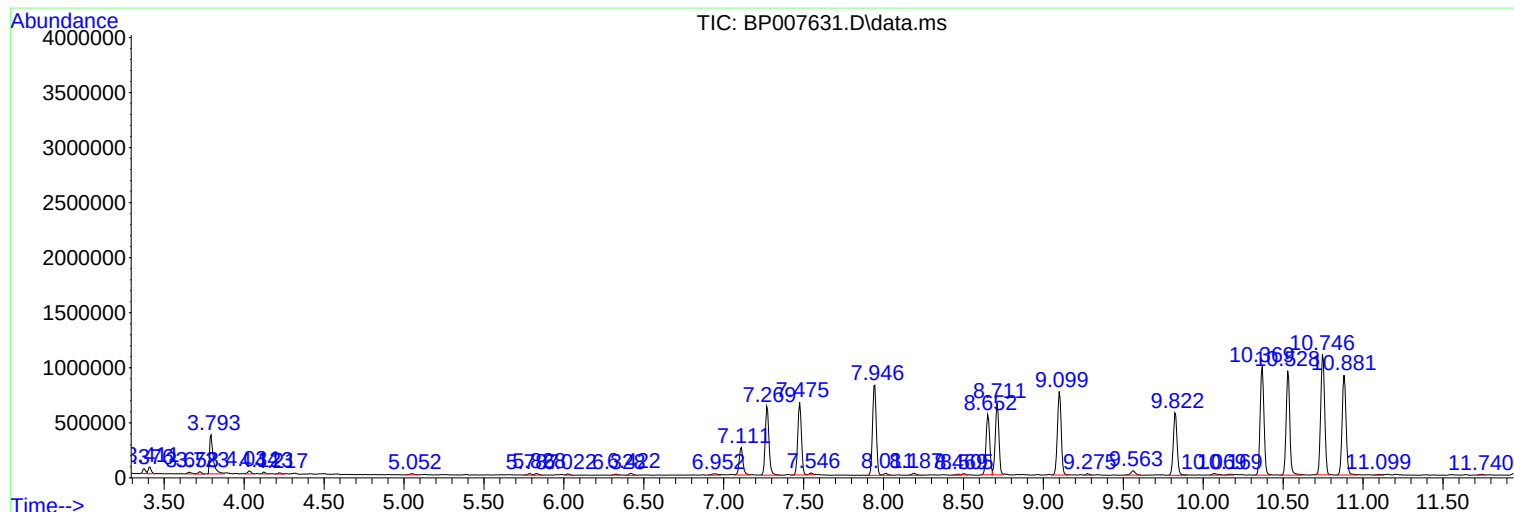
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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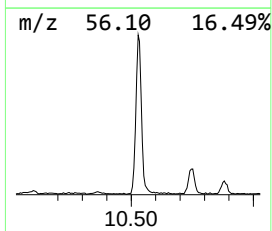
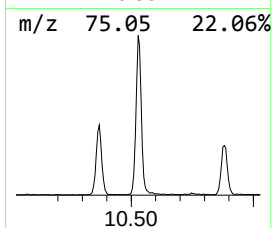
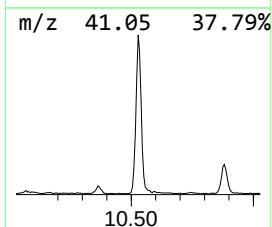
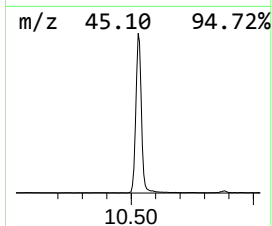
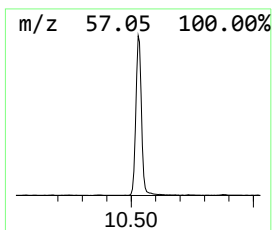
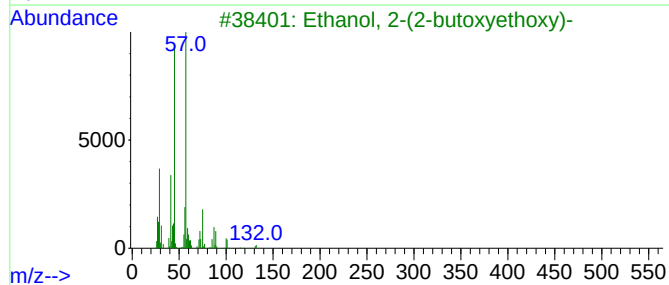
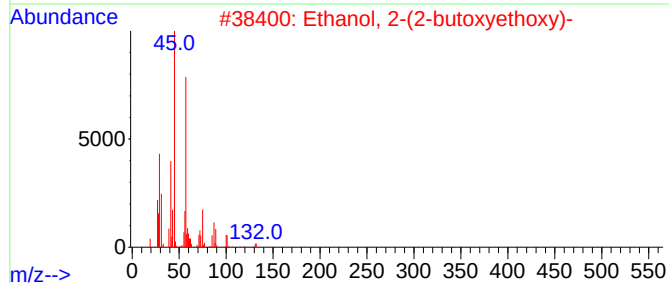
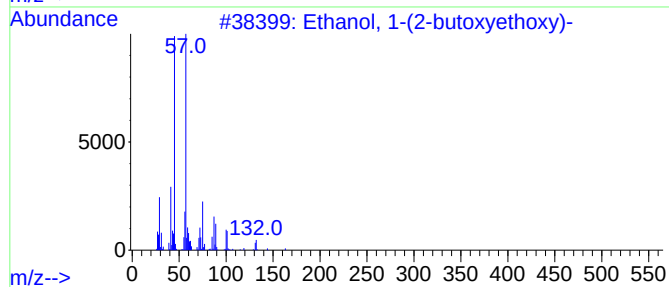
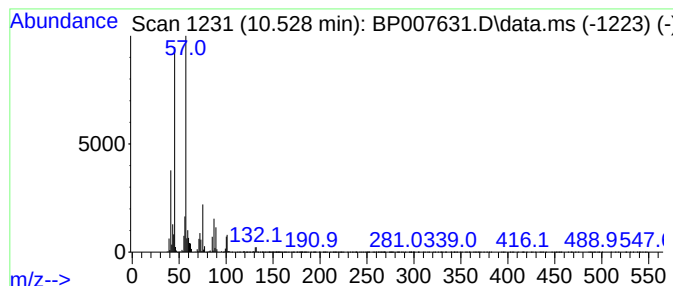
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.528	16.28 ng/ul	1503970	Naphthalene-d8	10.746

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	90
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72



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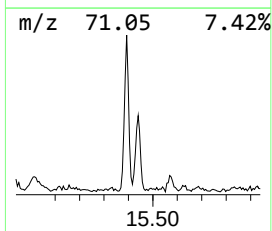
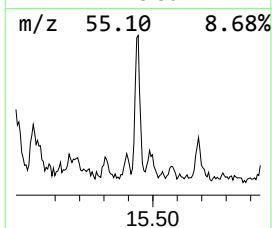
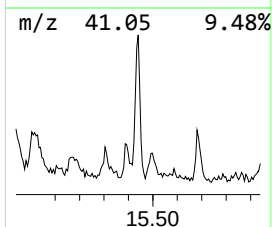
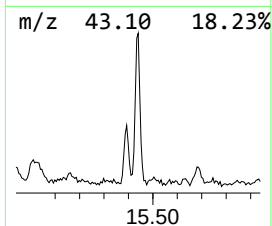
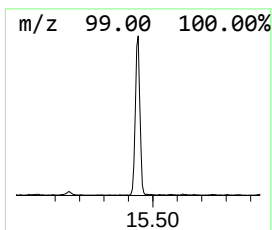
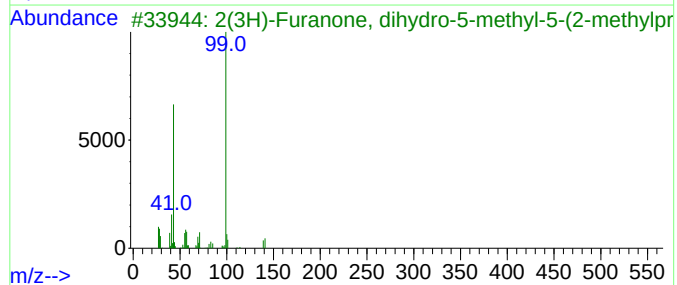
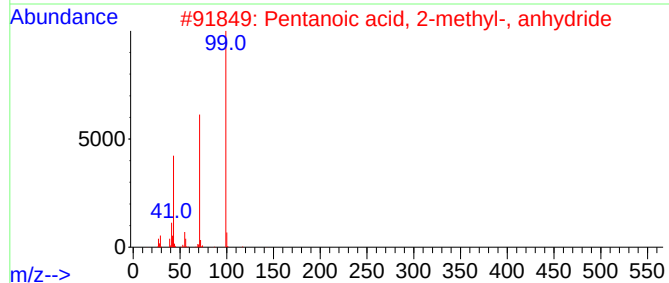
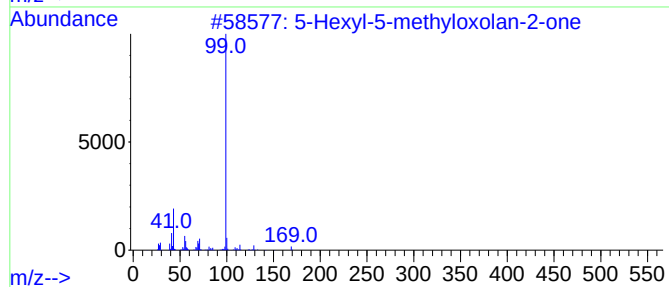
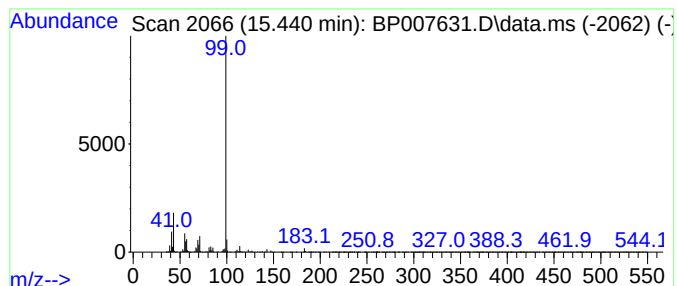
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 5-Hexyl-5-methyloxolan-2-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	2.01 ng/ul	283752	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Hexyl-5-methyloxolan-2-one	184	C11H20O2	007011-83-8	72
2			Pentanoic acid, 2-methyl-, anhyd...	214	C12H22O3	063169-61-9	64
3			2(3H)-Furanone, dihydro-5-methyl...	156	C9H16O2	010200-21-2	64
4			2-Pyrazoline-1-carboxaldehyde, 5...	184	C9H16N2O2	103214-44-4	56
5			Succinimide	99	C4H5NO2	000123-56-8	50



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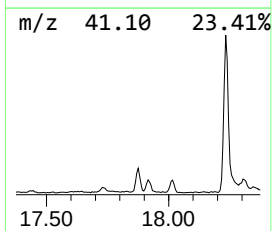
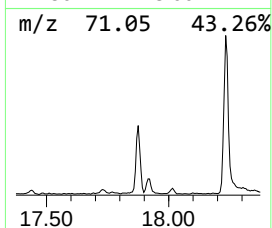
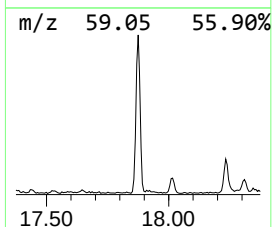
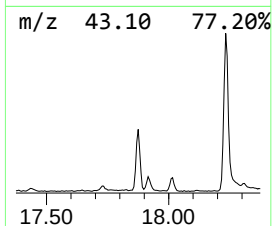
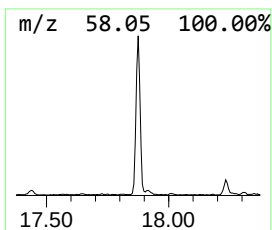
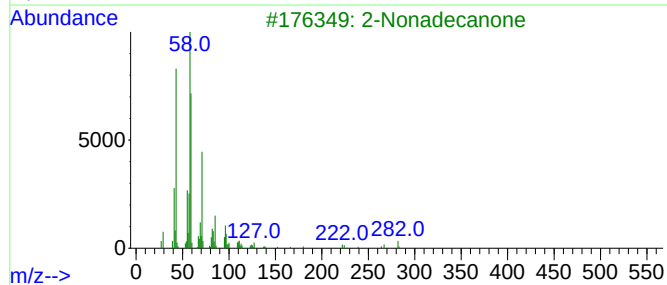
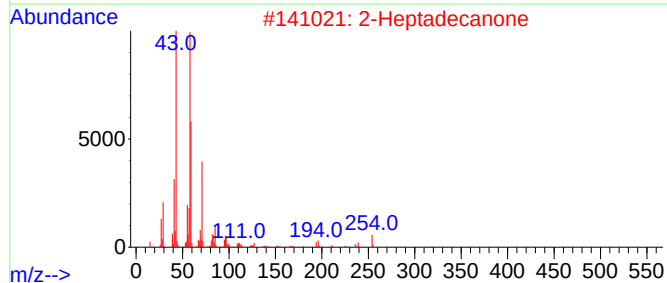
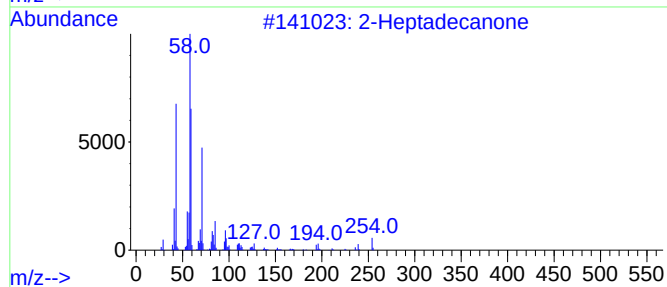
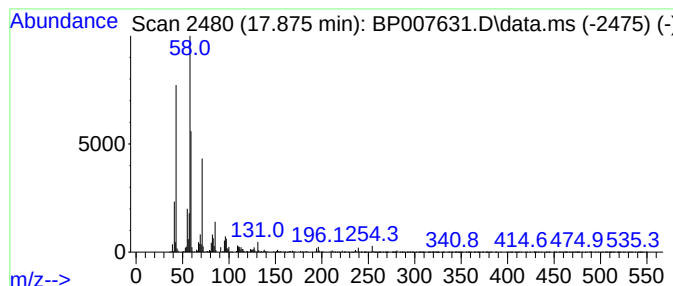
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Heptadecanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.875	2.65 ng/ul	405457	Phenanthrene-d10	17.334

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptadecanone			254	C17H34O	002922-51-2	97
2	2-Heptadecanone			254	C17H34O	002922-51-2	93
3	2-Nonadecanone			282	C19H38O	000629-66-3	90
4	2-Heptadecanone			254	C17H34O	002922-51-2	87
5	2-Pentadecanone			226	C15H30O	002345-28-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007631.D
Acq On : 23 Oct 2021 15:33
Operator : CG/JU
Sample : M4262-16
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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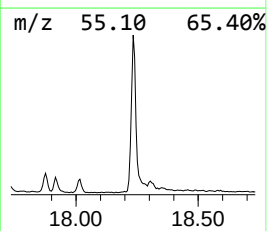
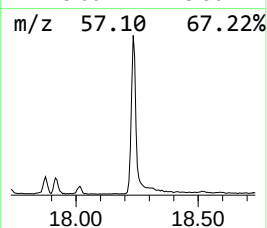
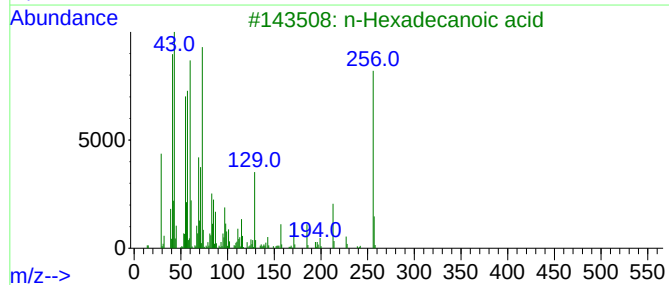
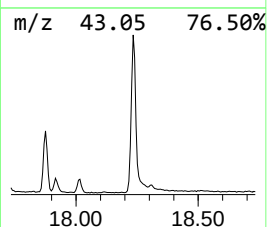
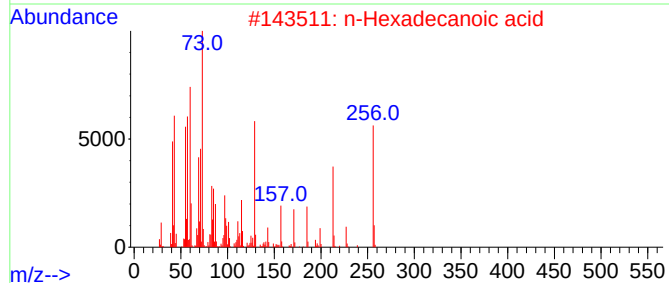
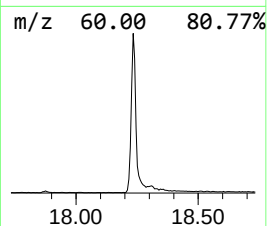
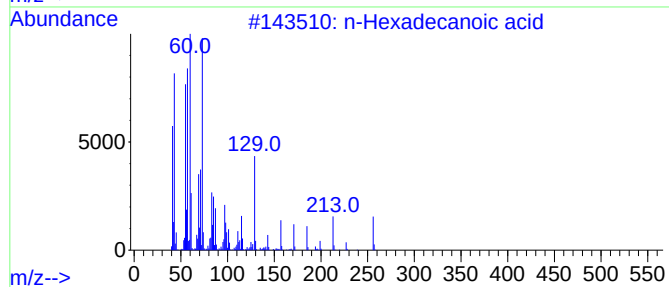
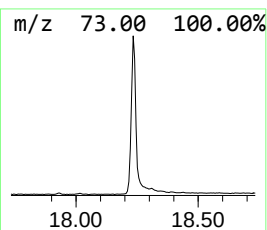
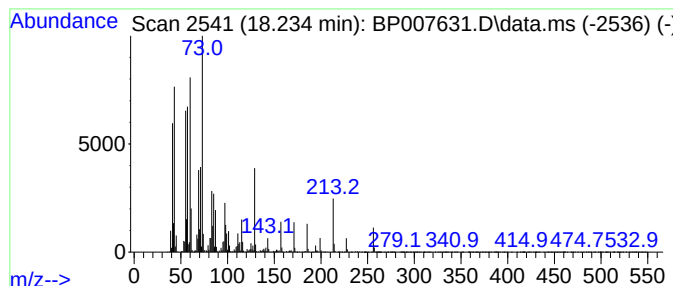
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.234	17.36 ng/ul	2652260	Phenanthrene-d10	17.334

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007631.D
Acq On : 23 Oct 2021 15:33
Operator : CG/JU
Sample : M4262-16
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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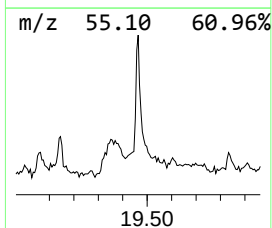
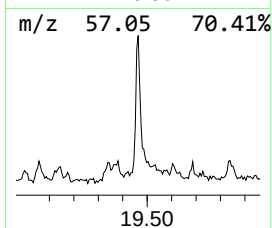
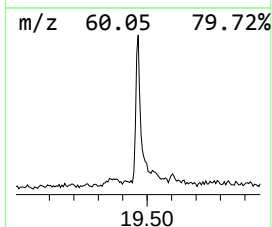
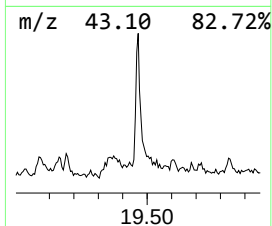
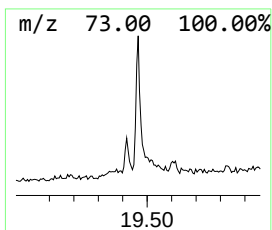
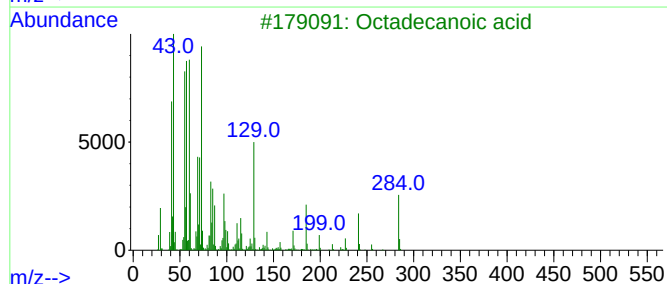
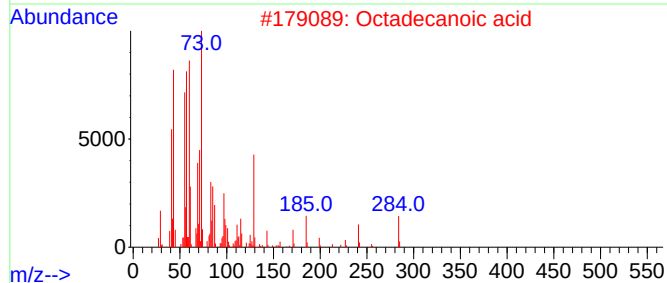
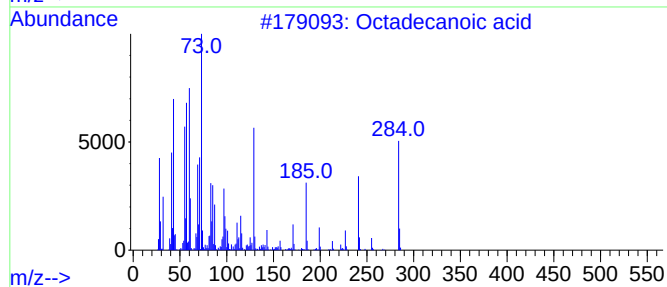
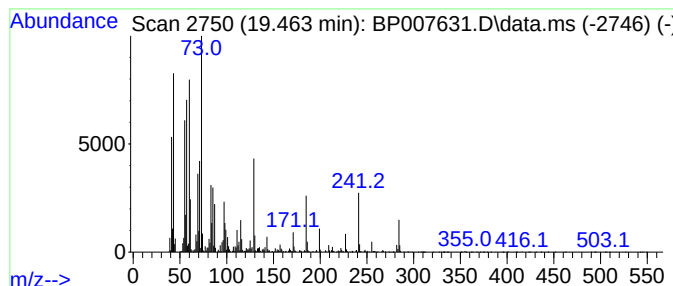
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	2.17 ng/ul	388204	Chrysene-d12	21.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	99
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007631.D
Acq On : 23 Oct 2021 15:33
Operator : CG/JU
Sample : M4262-16
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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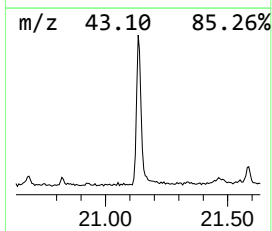
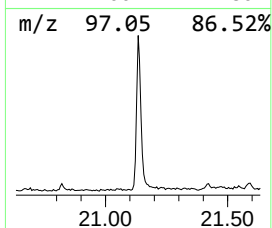
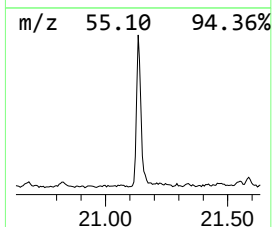
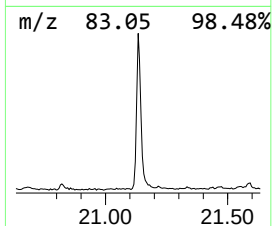
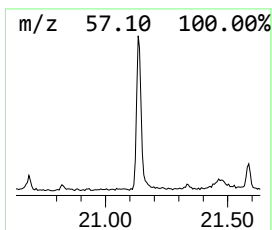
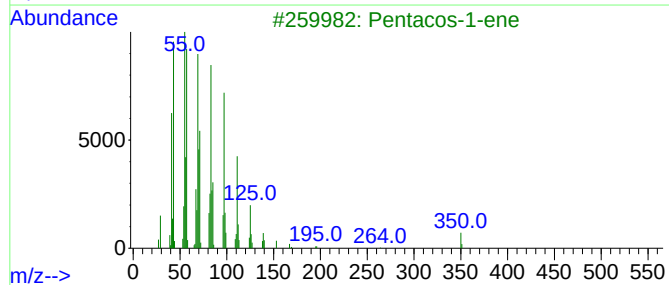
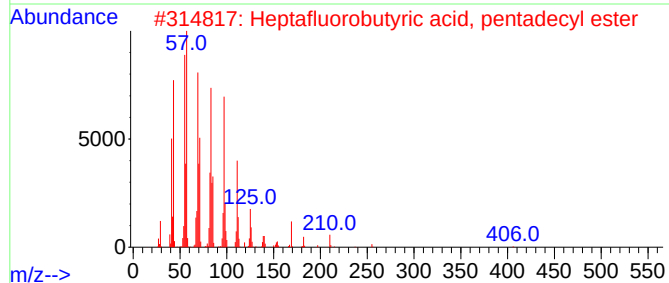
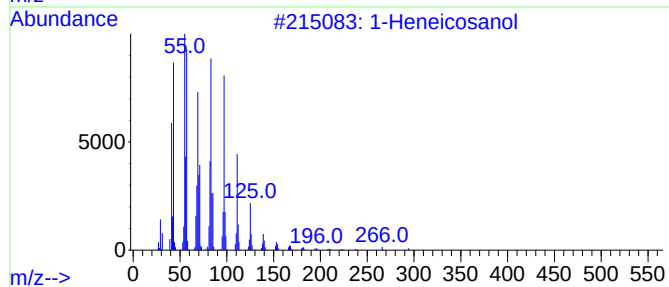
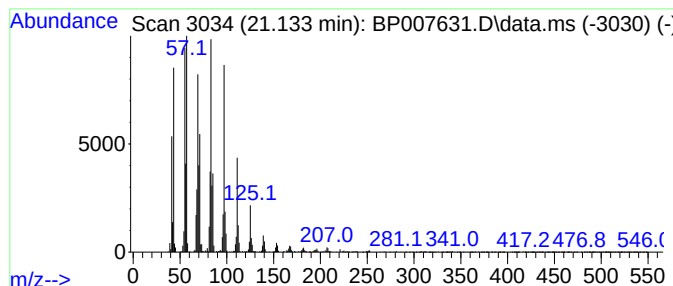
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	6.02 ng/ul	1076850	Chrysene-d12	21.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			1-Tetracosene	336	C24H48	010192-32-2	94
5			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007631.D
Acq On : 23 Oct 2021 15:33
Operator : CG/JU
Sample : M4262-16
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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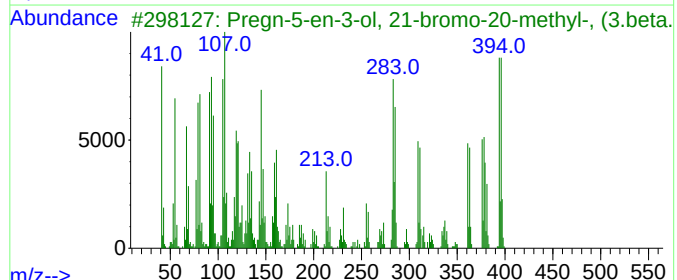
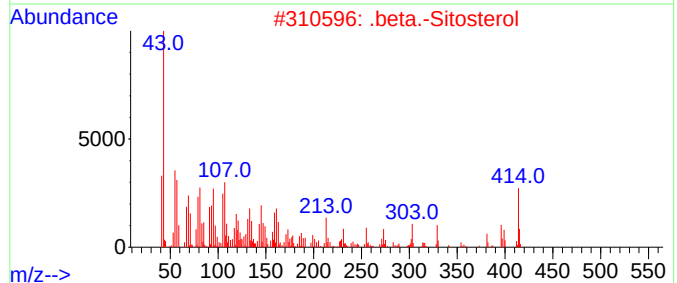
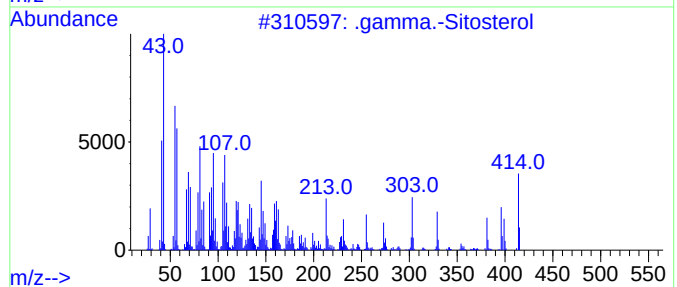
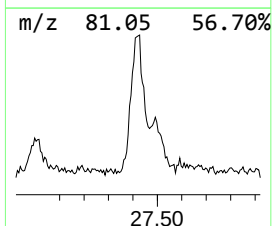
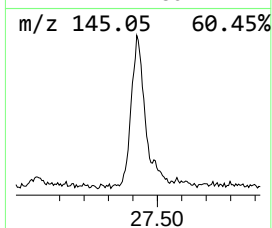
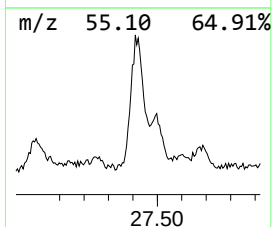
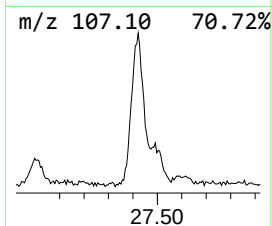
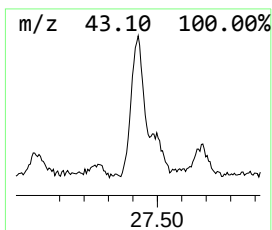
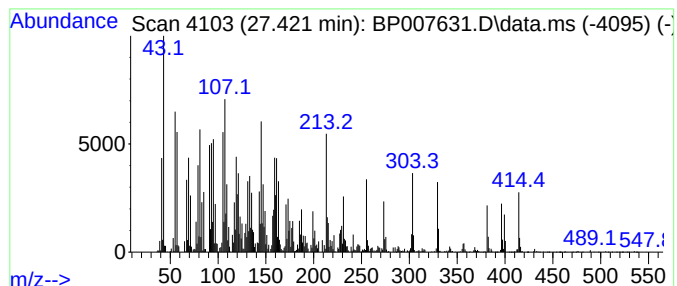
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 .gamma.-Sitosterol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.421	9.75 ng/ul	1799690	Perylene-d12	23.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	92
3			Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	78
4			5-Cholestene-3-ol, 24-methyl-	400	C28H48O	290299-12-6	49
5			.gamma.-Sitosterol	414	C29H50O	000083-47-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007631.D
Acq On : 23 Oct 2021 15:33
Operator : CG/JU
Sample : M4262-16
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.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
Ethanol, 1-(2-b...	10.528	16.3	ng/ul	1503970	2	10.746	1848010	20.0
5-Hexyl-5-methy...	15.440	2.0	ng/ul	283752	3	14.581	2826080	20.0
2-Heptadecanone	17.875	2.6	ng/ul	405457	4	17.334	3056340	20.0
n-Hexadecanoic ...	18.234	17.4	ng/ul	2652260	4	17.334	3056340	20.0
Octadecanoic acid	19.463	2.2	ng/ul	388204	5	21.422	3579470	20.0
1-Heneicosanol	21.133	6.0	ng/ul	1076850	5	21.422	3579470	20.0
.gamma.-Sitosterol	27.421	9.8	ng/ul	1799690	6	23.869	3690150	20.0