

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007499.D
 Acq On : 19 Oct 2021 18:12
 Operator : CG/JU
 Sample : M4196-03
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGA2

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: BP007499.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	12	15	25	rVB	44617	60088	1.43%	0.128%
2	3.652	59	62	67	rBV	10740	15021	0.36%	0.032%
3	3.799	84	87	100	rBV	196477	313400	7.44%	0.666%
4	4.040	125	128	133	rVB	13101	17224	0.41%	0.037%
5	4.134	140	144	149	rVB2	10678	14743	0.35%	0.031%
6	4.228	155	160	164	rBV7	7300	12709	0.30%	0.027%
7	4.458	197	199	203	rVB5	4773	5495	0.13%	0.012%
8	4.511	203	208	213	rBV9	3922	8224	0.20%	0.017%
9	5.058	296	301	310	rVB3	15560	24487	0.58%	0.052%
10	5.652	398	402	406	rVB7	3754	5997	0.14%	0.013%
11	6.040	465	468	472	rVB4	4905	5938	0.14%	0.013%
12	6.328	515	517	524	rVB6	6235	8259	0.20%	0.018%
13	6.940	617	621	623	rBV3	4293	5368	0.13%	0.011%
14	7.111	643	650	659	rBV	720890	1140053	27.07%	2.424%
15	7.281	672	679	688	rVB	569266	882590	20.96%	1.877%
16	7.481	705	713	723	rBV	709634	1144565	27.18%	2.434%
17	7.558	723	726	730	rVV3	8403	13274	0.32%	0.028%
18	7.958	786	794	801	rBV	639593	1028895	24.43%	2.188%
19	8.028	801	806	811	rVB5	5761	9724	0.23%	0.021%
20	8.658	906	913	923	rBV	899035	1415858	33.62%	3.010%
21	9.111	980	990	1001	rBV	649086	1055314	25.06%	2.244%
22	9.287	1016	1020	1024	rVB7	7024	8617	0.20%	0.018%
23	9.581	1066	1070	1075	rVB5	7452	10744	0.26%	0.023%
24	9.834	1106	1113	1123	rBV	491767	801911	19.04%	1.705%
25	10.375	1198	1205	1217	rBV	851535	1424178	33.82%	3.028%
26	10.540	1227	1233	1248	rBV	430466	674216	16.01%	1.433%
27	10.763	1262	1271	1280	rVB	834242	1448691	34.40%	3.080%
28	10.887	1282	1292	1302	rBV	798767	1367613	32.48%	2.908%
29	11.558	1403	1406	1407	rBV3	3657	4399	0.10%	0.009%
30	12.281	1523	1529	1534	rBV5	10266	17491	0.42%	0.037%
31	12.346	1534	1540	1549	rVB2	18259	31653	0.75%	0.067%
32	12.563	1572	1577	1581	rBV8	2985	5719	0.14%	0.012%
33	12.963	1641	1645	1650	rBV6	6328	9444	0.22%	0.020%
34	13.099	1663	1668	1672	rBV8	2651	5239	0.12%	0.011%
35	13.210	1681	1687	1691	rBV3	10598	15756	0.37%	0.033%
36	13.499	1731	1736	1742	rBV4	12653	20823	0.49%	0.044%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007499.D
 Acq On : 19 Oct 2021 18:12
 Operator : CG/JU
 Sample : M4196-03
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGA2

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

37	13.563	1742	1747	1754	rVB6	5850	10302	0.24%	0.022%
38	13.793	1782	1786	1791	rBV7	10536	18320	0.44%	0.039%
39	13.999	1814	1821	1828	rBV	1449090	2009986	47.73%	4.274%
40	14.281	1860	1869	1878	rBV	1679802	2501687	59.41%	5.319%
41	14.510	1905	1908	1914	rVB8	3636	5143	0.12%	0.011%
42	14.593	1914	1922	1929	rBV	1308802	1968195	46.74%	4.185%
43	14.769	1944	1952	1968	rBV	664926	987313	23.45%	2.099%
44	15.004	1985	1992	1998	rBV	9506	15396	0.37%	0.033%
45	15.240	2027	2032	2034	rBV6	5312	6628	0.16%	0.014%
46	15.410	2054	2061	2065	rBV3	9796	15633	0.37%	0.033%
47	15.457	2065	2069	2073	rVV7	4176	5042	0.12%	0.011%
48	15.581	2083	2090	2101	rBV	2119456	3118076	74.05%	6.630%
49	15.693	2103	2109	2116	rBV	367043	504232	11.97%	1.072%
50	15.822	2126	2131	2137	rVB2	27521	43229	1.03%	0.092%
51	15.957	2150	2154	2157	rBV6	3060	4242	0.10%	0.009%
52	16.028	2161	2166	2173	rBV6	15725	24997	0.59%	0.053%
53	16.140	2182	2185	2193	rBV9	7187	14022	0.33%	0.030%
54	16.263	2199	2206	2213	rBV9	7424	19763	0.47%	0.042%
55	16.375	2220	2225	2227	rBV6	5808	8469	0.20%	0.018%
56	16.493	2239	2245	2250	rVB8	8149	14769	0.35%	0.031%
57	16.593	2260	2262	2266	rVB5	4924	4753	0.11%	0.010%
58	16.751	2285	2289	2293	rVB7	7669	10559	0.25%	0.022%
59	16.951	2319	2323	2328	rBV8	3269	5756	0.14%	0.012%
60	17.022	2331	2335	2337	rBV5	4162	5054	0.12%	0.011%
61	17.122	2344	2352	2357	rBV6	11353	22119	0.53%	0.047%
62	17.281	2376	2379	2382	rBV5	3463	4489	0.11%	0.010%
63	17.340	2382	2389	2397	rVV	1626556	2427006	57.64%	5.160%
64	17.440	2400	2406	2417	rVV	2366456	3421415	81.25%	7.274%
65	17.534	2418	2422	2426	rVV7	16878	29050	0.69%	0.062%
66	17.575	2426	2429	2436	rVB5	12150	21467	0.51%	0.046%
67	18.028	2501	2506	2512	rBV	108372	129895	3.08%	0.276%
68	18.234	2536	2541	2550	rBV	163987	242960	5.77%	0.517%
69	18.410	2569	2571	2575	rBV4	3886	4350	0.10%	0.009%
70	18.534	2590	2592	2595	rBV4	4932	5025	0.12%	0.011%
71	19.157	2694	2698	2704	rVB2	84927	96520	2.29%	0.205%
72	19.251	2710	2714	2717	rBV3	32590	40336	0.96%	0.086%
73	19.287	2718	2720	2722	rVV3	12228	13026	0.31%	0.028%
74	19.322	2722	2726	2729	rVV	44056	63707	1.51%	0.135%
75	19.357	2729	2732	2740	rVB7	27946	56775	1.35%	0.121%
76	19.469	2747	2751	2756	rBV	82763	110118	2.62%	0.234%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007499.D
 Acq On : 19 Oct 2021 18:12
 Operator : CG/JU
 Sample : M4196-03
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGA2

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

77	19.598	2770	2773	2775	rBV3	21526	31009	0.74%	0.066%
78	19.675	2780	2786	2792	rBV	3075731	4210670	99.99%	8.953%
79	20.581	2936	2940	2946	rBV	128066	159928	3.80%	0.340%
80	21.145	3032	3036	3041	rBV2	364977	493440	11.72%	1.049%
81	21.428	3079	3084	3093	rBV	2307105	2961387	70.33%	6.296%
82	22.575	3274	3279	3294	rBV3	475734	875364	20.79%	1.861%
83	23.722	3467	3474	3485	rVB2	2050848	4210915	100.00%	8.953%
84	23.880	3495	3501	3511	rVB2	1464541	3046963	72.36%	6.478%

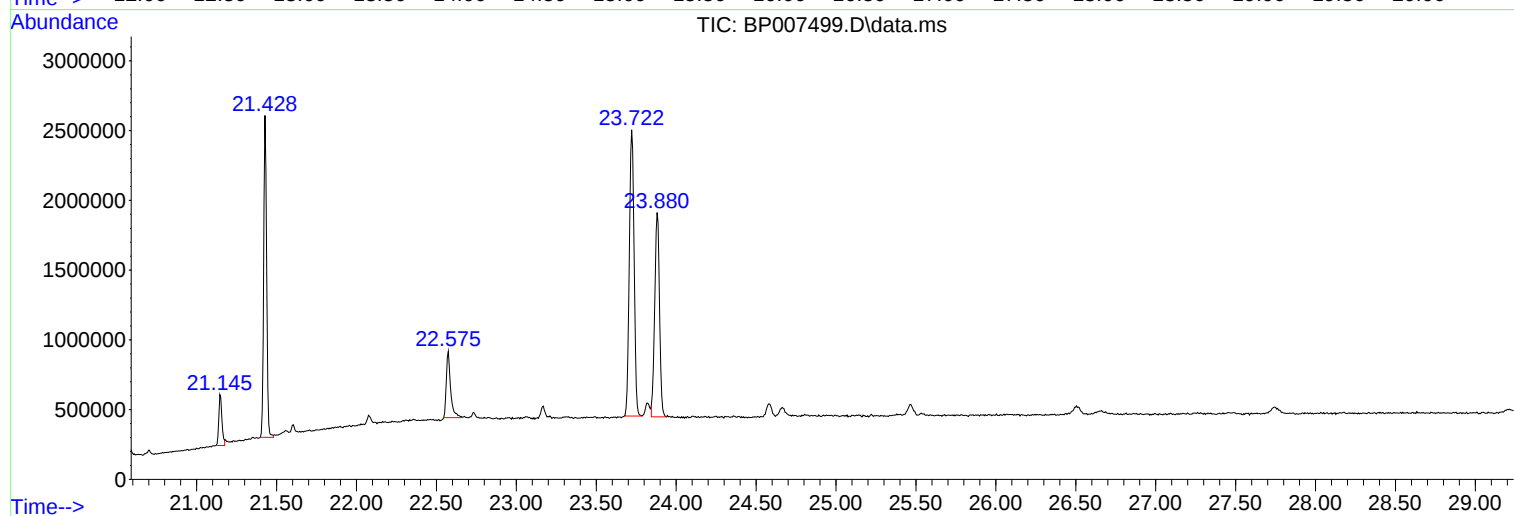
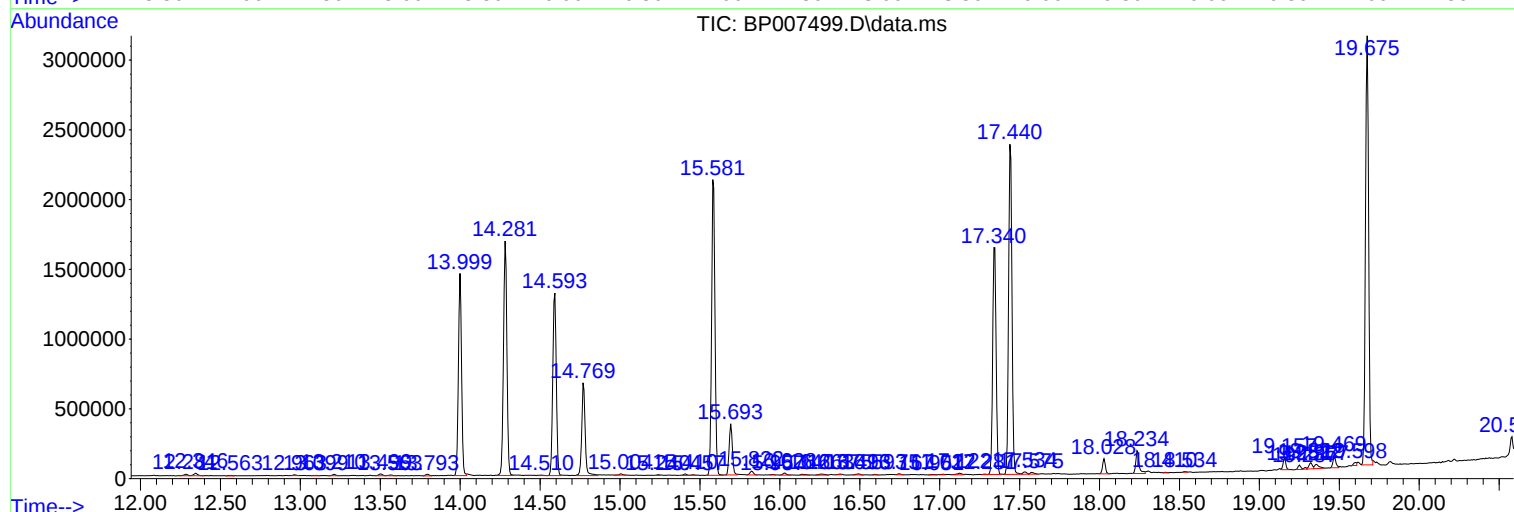
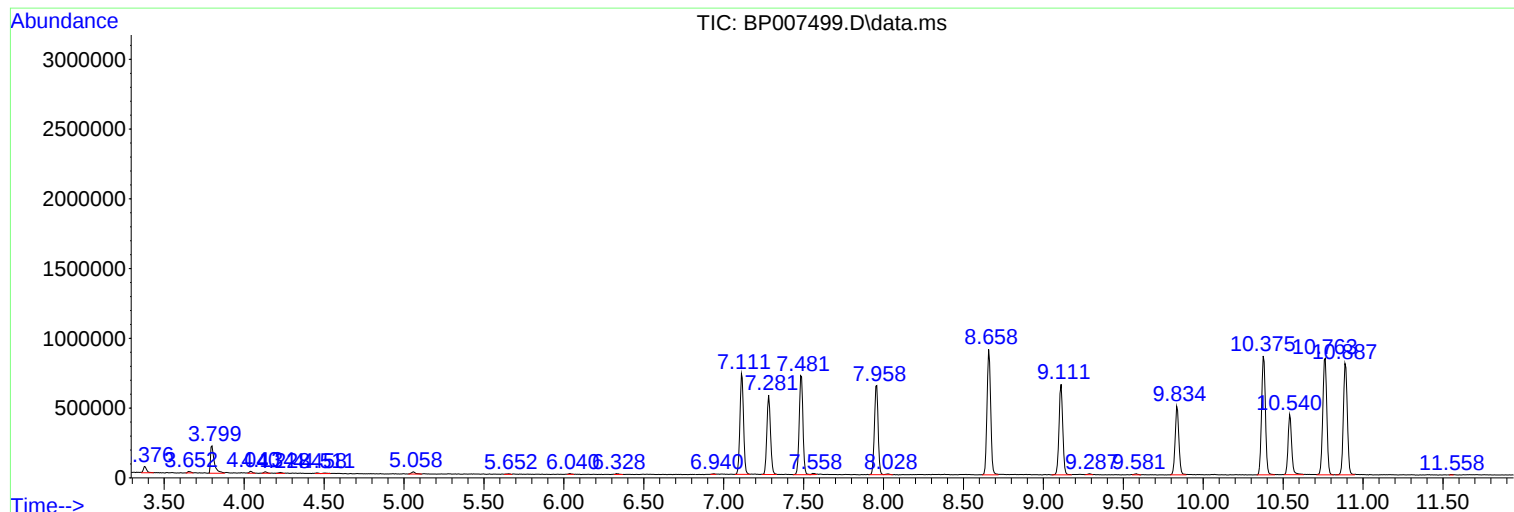
Sum of corrected areas: 47033220

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

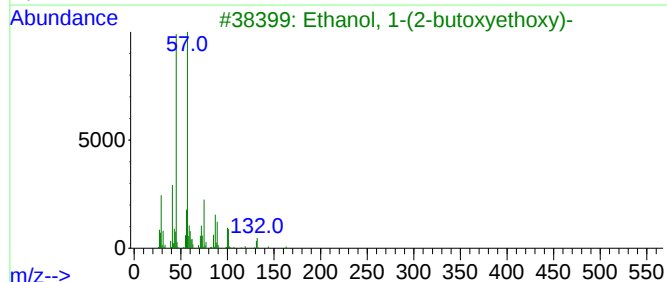
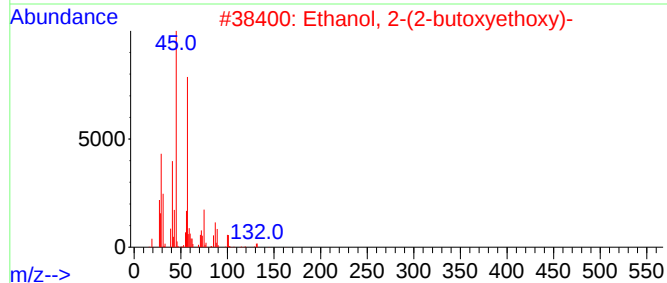
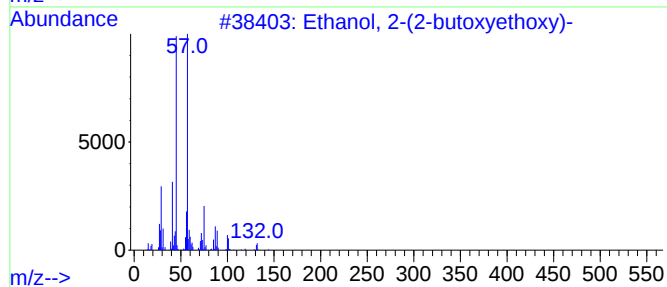
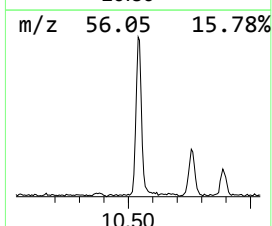
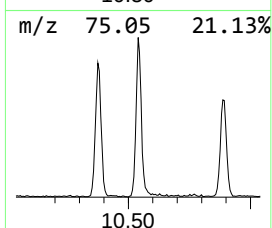
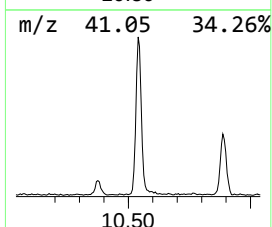
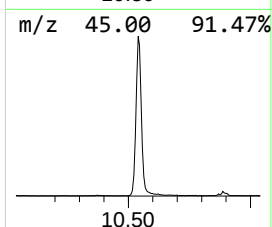
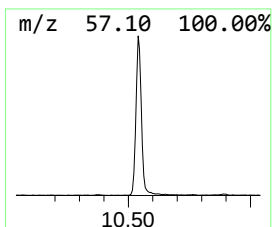
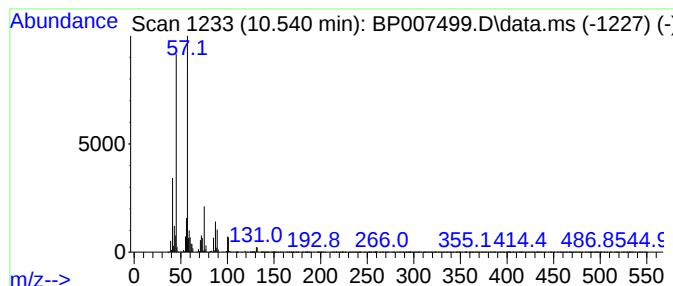
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	9.31 ng/ul	674216	Naphthalene-d8	10.763

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

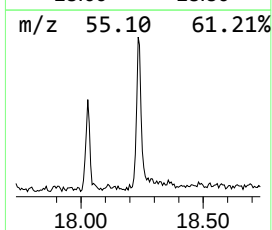
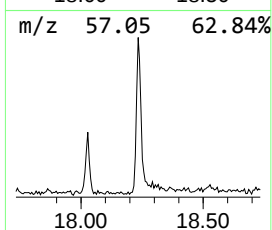
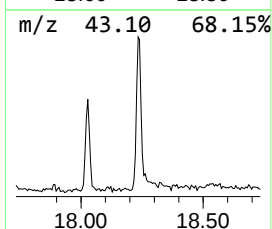
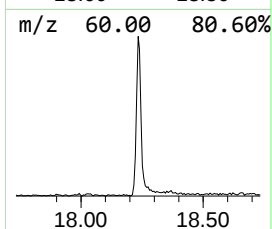
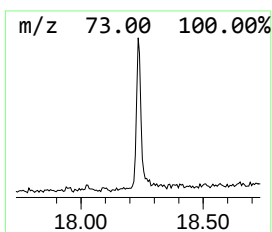
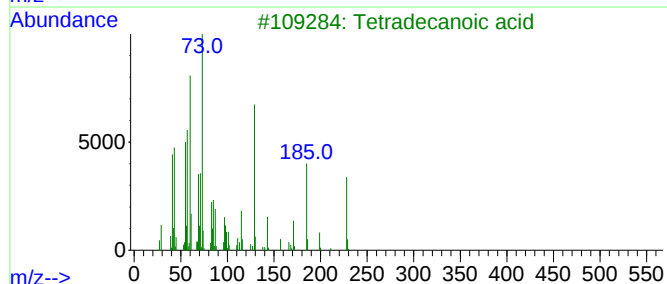
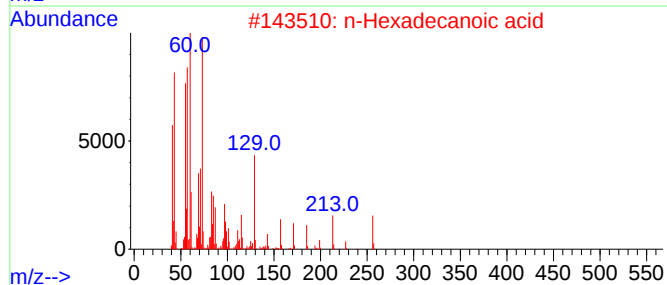
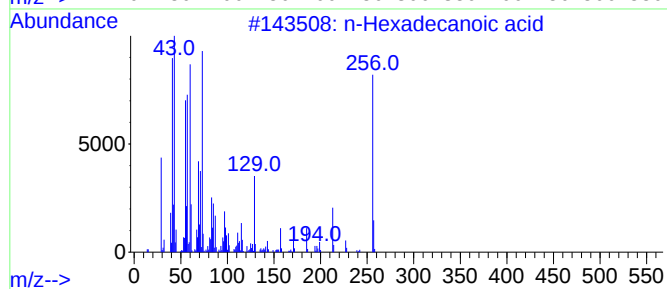
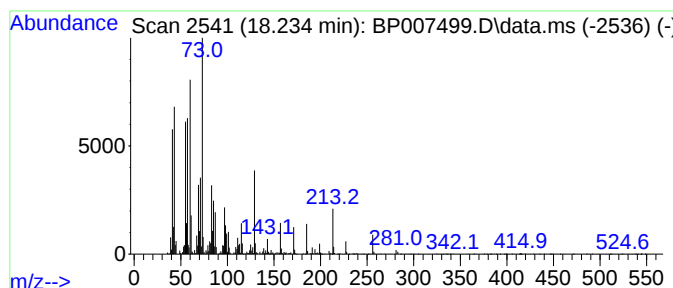
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.234	2.00 ng/ul	242960	Phenanthrene-d10	17.345

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	95
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

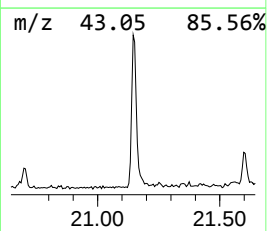
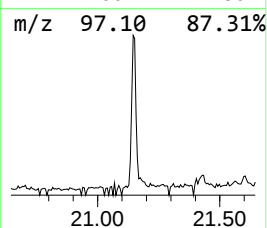
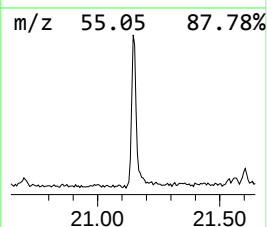
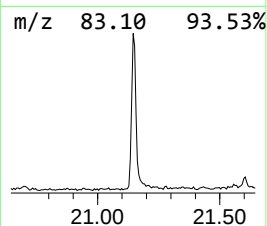
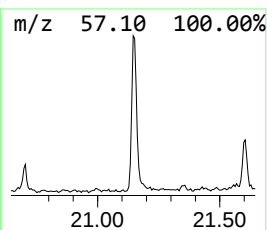
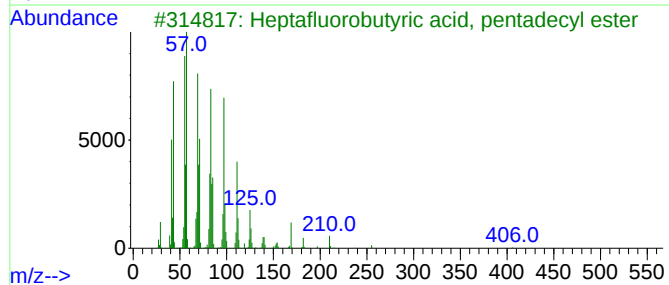
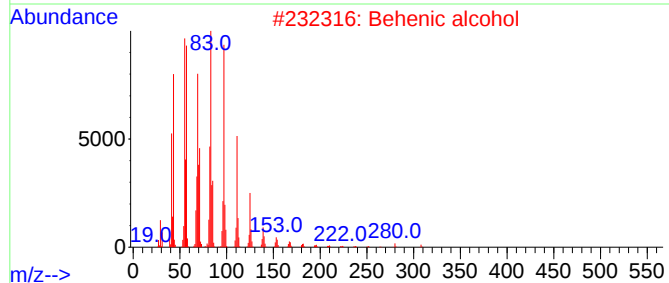
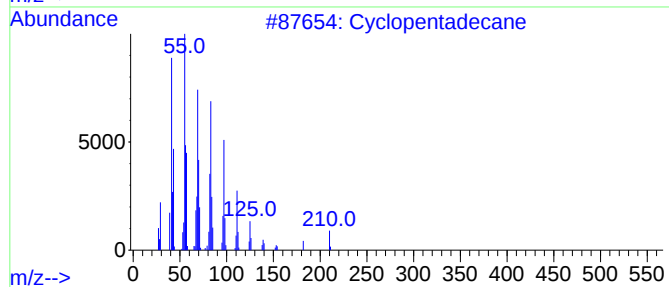
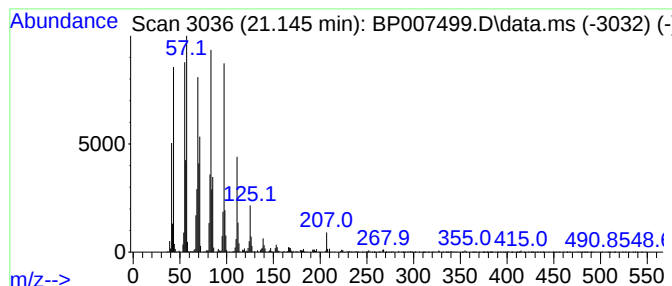
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic21.145 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	3.33 ng/ul	493440	Chrysene-d12	21.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	94
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
5			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

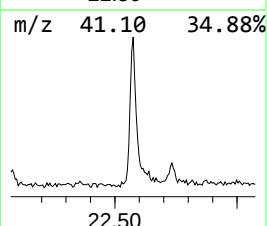
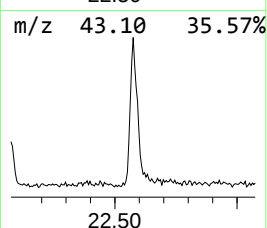
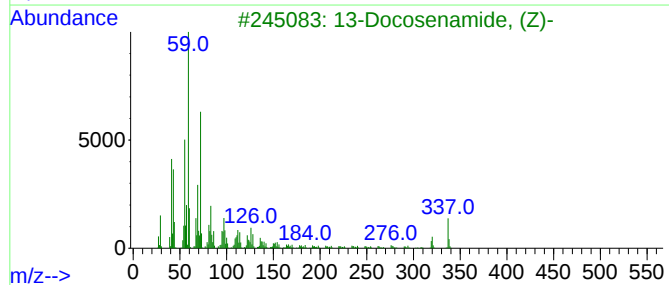
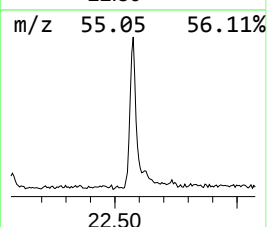
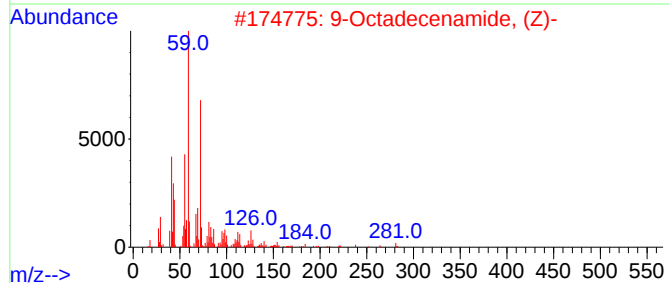
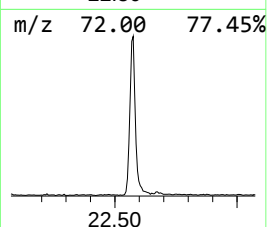
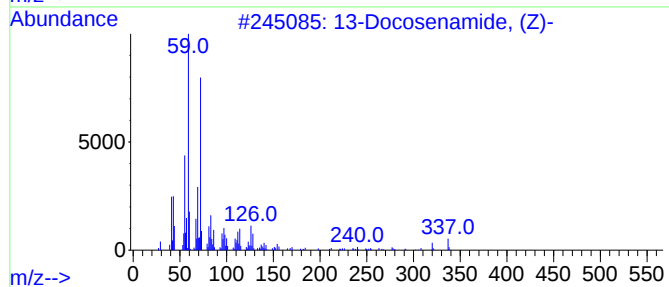
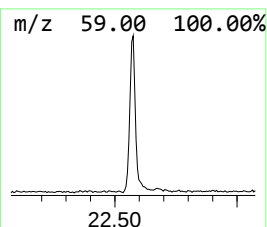
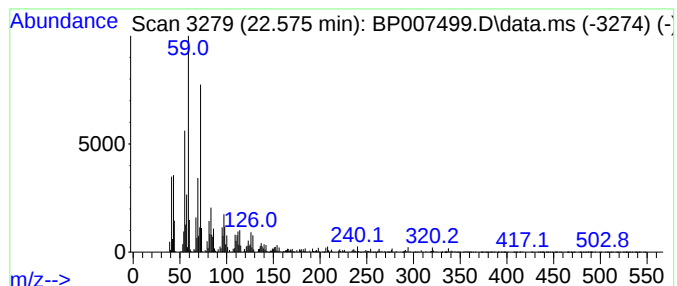
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 13-Docosenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.575	5.91 ng/ul	875364	Chrysene-d12	21.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	98
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	83
4			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
5			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007499.D
Acq On : 19 Oct 2021 18:12
Operator : CG/JU
Sample : M4196-03
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA2

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.540	9.3	ng/ul	674216	2	10.763	1448690	20.0
n-Hexadecanoic ...	18.234	2.0	ng/ul	242960	4	17.345	2427010	20.0
(DEL) Alkane: C...	21.145	3.3	ng/ul	493440	5	21.428	2961390	20.0
13-Docosenamide...	22.575	5.9	ng/ul	875364	5	21.428	2961390	20.0