

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007494.D
 Acq On : 19 Oct 2021 15:21
 Operator : CG/JU
 Sample : M4196-05
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGA4

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: BP007494.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	11	15	22	rVB	37375	54876	1.01%	0.094%
2	3.652	59	62	67	rBV	19628	30090	0.55%	0.051%
3	3.799	84	87	98	rBV	176888	254405	4.68%	0.434%
4	4.040	125	128	133	rVB2	13631	16663	0.31%	0.028%
5	4.134	140	144	147	rBV	10268	14420	0.27%	0.025%
6	5.058	297	301	306	rVB2	12109	17818	0.33%	0.030%
7	5.652	399	402	407	rBV6	3637	5436	0.10%	0.009%
8	6.016	459	464	466	rBV6	10378	18234	0.34%	0.031%
9	6.334	514	518	523	rVB7	7177	12135	0.22%	0.021%
10	6.946	618	622	630	rVB10	5895	13653	0.25%	0.023%
11	7.116	644	651	660	rBV	913355	1412104	26.00%	2.410%
12	7.281	673	679	690	rVB	662191	1065738	19.62%	1.819%
13	7.487	705	714	722	rBV	880308	1406990	25.91%	2.402%
14	7.563	724	727	730	rBV4	5375	6965	0.13%	0.012%
15	7.957	788	794	801	rVV	675319	1055008	19.43%	1.801%
16	8.028	802	806	810	rVB3	6899	9853	0.18%	0.017%
17	8.663	904	914	924	rBV	1128516	1828682	33.67%	3.121%
18	9.110	983	990	997	rBV	820147	1339635	24.67%	2.287%
19	9.287	1017	1020	1025	rVB7	5924	8897	0.16%	0.015%
20	9.581	1063	1070	1077	rBV10	6732	14525	0.27%	0.025%
21	9.840	1105	1114	1123	rBV	636408	1061597	19.55%	1.812%
22	10.381	1198	1206	1214	rBV	1092965	1843817	33.95%	3.147%
23	10.546	1227	1234	1245	rBV	210017	345583	6.36%	0.590%
24	10.763	1263	1271	1280	rBV	876938	1502970	27.67%	2.565%
25	10.893	1284	1293	1303	rBV	1074042	1779093	32.76%	3.037%
26	11.128	1328	1333	1338	rBV9	2451	7128	0.13%	0.012%
27	11.557	1401	1406	1415	rVB10	8525	16390	0.30%	0.028%
28	12.287	1526	1530	1536	rVV6	4819	8695	0.16%	0.015%
29	12.345	1536	1540	1547	rVB3	18899	30701	0.57%	0.052%
30	12.557	1573	1576	1581	rBV5	3922	6480	0.12%	0.011%
31	12.963	1640	1645	1648	rBV6	5802	8232	0.15%	0.014%
32	13.216	1684	1688	1692	rVB6	7356	10199	0.19%	0.017%
33	13.439	1722	1726	1731	rVB7	5672	8691	0.16%	0.015%
34	13.498	1731	1736	1743	rBV9	8477	17035	0.31%	0.029%
35	13.563	1743	1747	1752	rVB6	6309	11517	0.21%	0.020%
36	13.934	1803	1810	1815	rBV8	16182	32862	0.61%	0.056%

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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 >

Peak separation: 5

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

37	14.004	1815	1822	1836	rVB	1797720	2592327	47.73%	4.425%
38	14.175	1848	1851	1856	rVB7	5486	7078	0.13%	0.012%
39	14.286	1858	1870	1878	rBV	2111003	3232330	59.52%	5.517%
40	14.592	1915	1922	1935	rVB	1361167	2038107	37.53%	3.479%
41	14.775	1945	1953	1966	rBV	910663	1304716	24.02%	2.227%
42	15.004	1989	1992	1997	rVB5	17472	24554	0.45%	0.042%
43	15.263	2029	2036	2042	rBV7	7713	17056	0.31%	0.029%
44	15.404	2055	2060	2065	rBV5	7841	14534	0.27%	0.025%
45	15.463	2066	2070	2074	rVB5	4063	6818	0.13%	0.012%
46	15.586	2084	2091	2102	rVB	2683308	3936959	72.49%	6.720%
47	15.692	2103	2109	2120	rBV	491382	711357	13.10%	1.214%
48	15.828	2127	2132	2138	rVB2	33183	49925	0.92%	0.085%
49	15.957	2150	2154	2155	rBV4	5037	6211	0.11%	0.011%
50	16.033	2163	2167	2172	rVB8	5612	9109	0.17%	0.016%
51	16.151	2181	2187	2193	rBV8	4689	9668	0.18%	0.017%
52	16.269	2204	2207	2213	rVV7	6913	10450	0.19%	0.018%
53	16.386	2221	2227	2231	rVB9	8364	16743	0.31%	0.029%
54	16.751	2286	2289	2293	rVB6	8151	9957	0.18%	0.017%
55	17.022	2331	2335	2337	rBV5	5678	6911	0.13%	0.012%
56	17.133	2345	2354	2357	rBV4	12516	25672	0.47%	0.044%
57	17.345	2383	2390	2399	rVB	1737326	2483820	45.73%	4.240%
58	17.445	2400	2407	2417	rVV	2994657	4239759	78.07%	7.237%
59	17.539	2418	2423	2427	rVV5	27771	46133	0.85%	0.079%
60	17.580	2427	2430	2436	rVB5	21684	32223	0.59%	0.055%
61	17.886	2478	2482	2485	rBV5	4769	8577	0.16%	0.015%
62	18.027	2502	2506	2512	rBV	32521	43757	0.81%	0.075%
63	18.239	2537	2542	2551	rBV	308193	430459	7.93%	0.735%
64	19.098	2684	2688	2697	rBV	144753	214397	3.95%	0.366%
65	19.251	2711	2714	2719	rBV2	37565	49773	0.92%	0.085%
66	19.327	2722	2727	2729	rBV3	119247	157558	2.90%	0.269%
67	19.357	2729	2732	2744	rVV2	145561	306476	5.64%	0.523%
68	19.469	2747	2751	2762	rVV	255595	386816	7.12%	0.660%
69	19.604	2770	2774	2775	rBV2	77966	96778	1.78%	0.165%
70	19.674	2781	2786	2791	rBV	3866871	5175601	95.30%	8.834%
71	19.821	2808	2811	2818	rVB3	55929	78978	1.45%	0.135%
72	20.392	2903	2908	2920	rBV	603900	1031913	19.00%	1.761%
73	20.580	2934	2940	2948	rBV	565594	730508	13.45%	1.247%
74	21.151	3033	3037	3042	rBV	371589	497105	9.15%	0.848%
75	21.427	3079	3084	3091	rBV	2159855	2931763	53.98%	5.004%
76	22.357	3239	3242	3248	rVB2	237834	306761	5.65%	0.524%

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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

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Peak separation: 5

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Title : SVOA CALIBRATION

77	22.574	3274	3279	3291	rBV3	664770	1236790	22.77%	2.111%
78	23.168	3377	3380	3386	rVB2	142762	215185	3.96%	0.367%
79	23.727	3468	3475	3485	rVB2	2659874	5431057	100.00%	9.270%
80	23.880	3495	3501	3512	rVB2	1444498	3147251	57.95%	5.372%

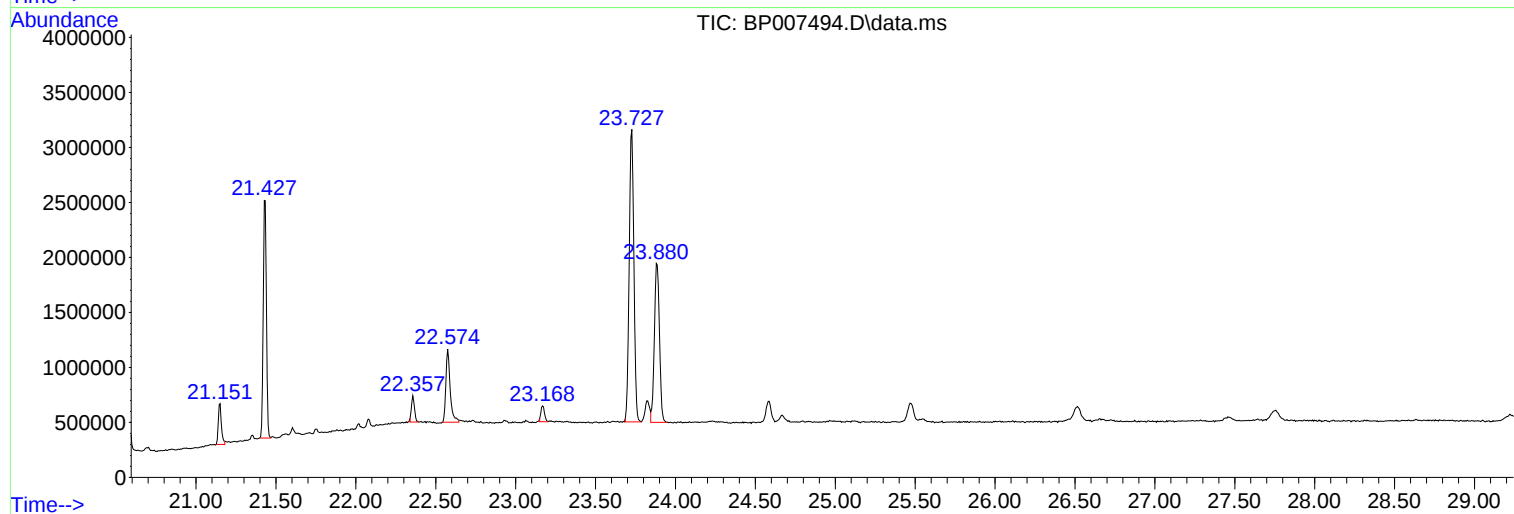
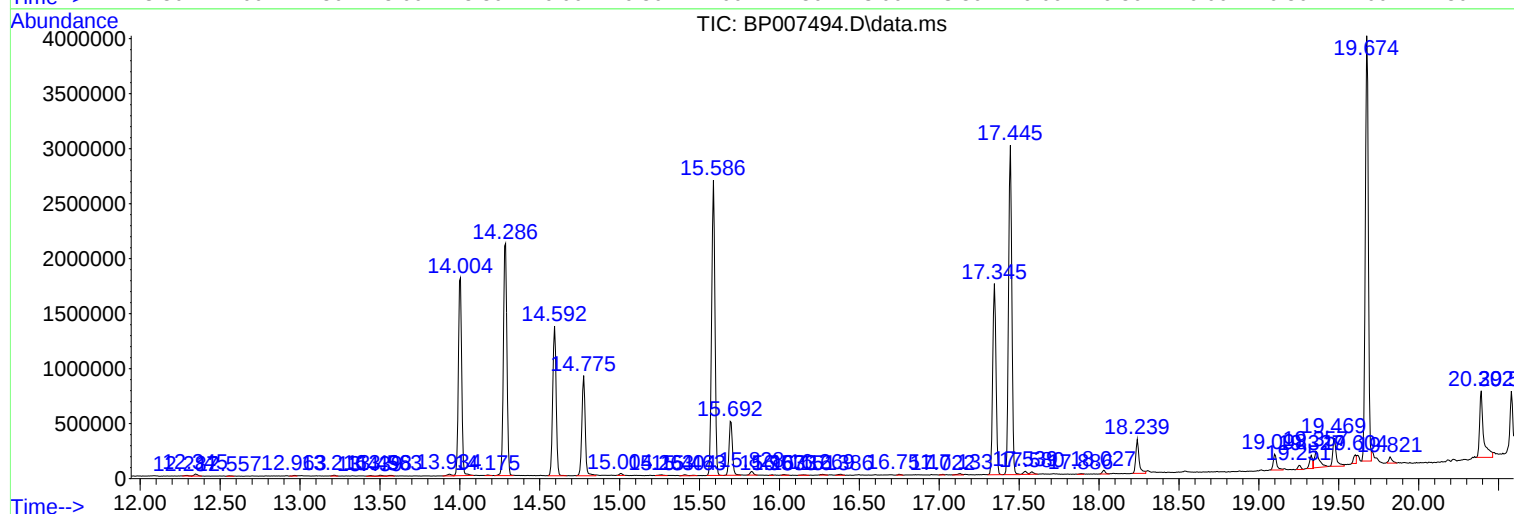
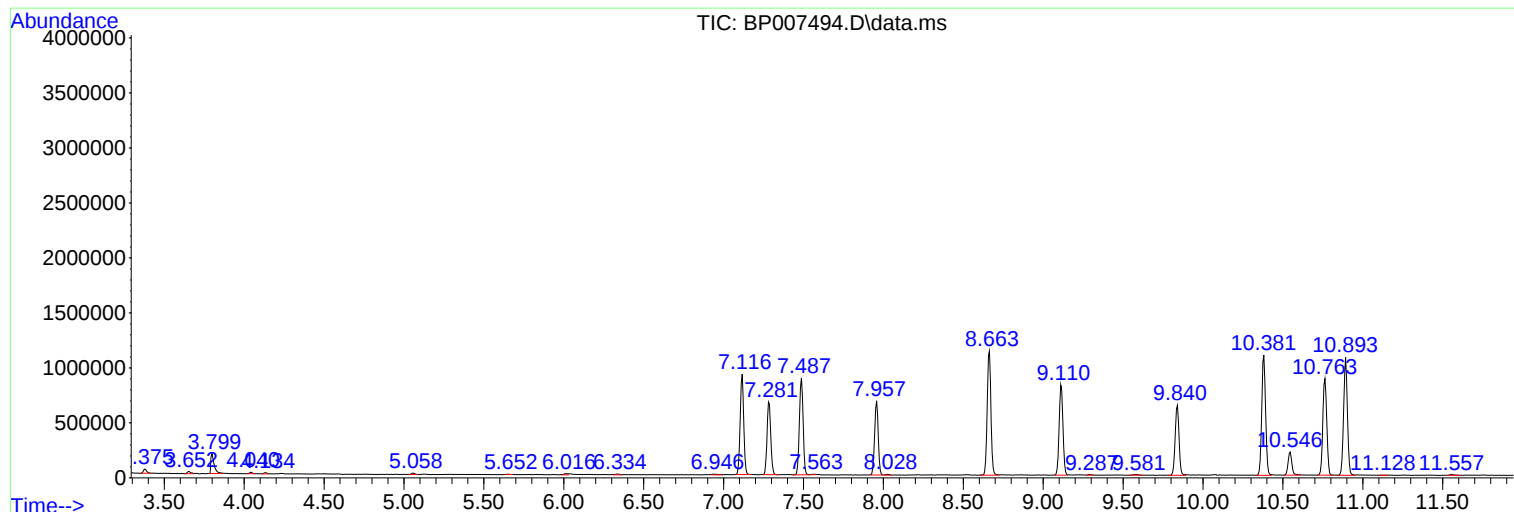
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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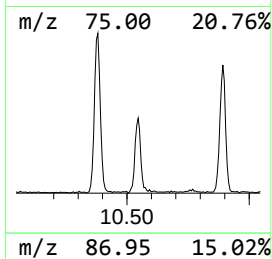
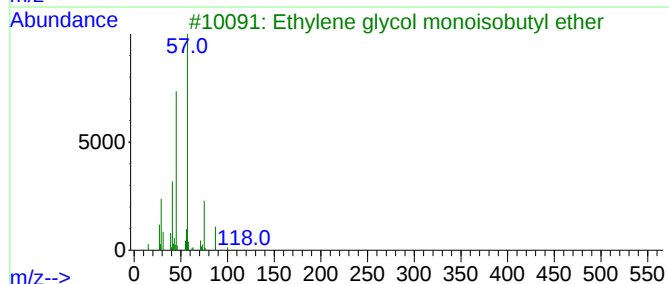
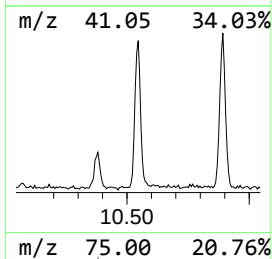
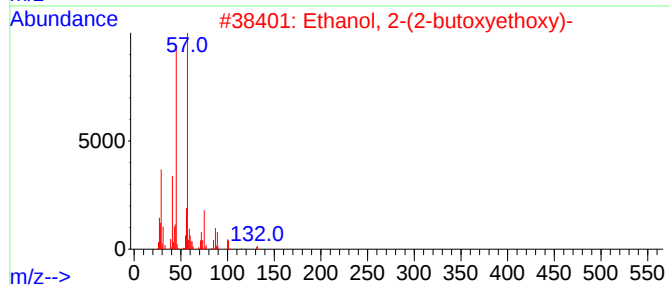
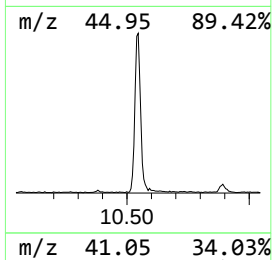
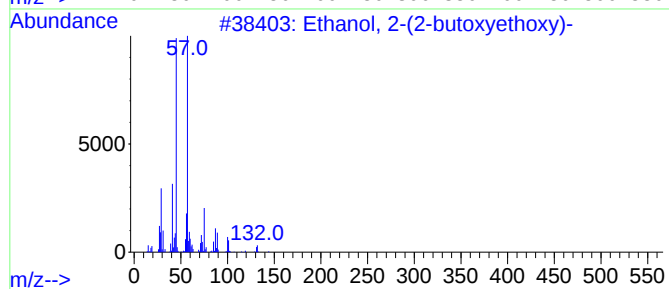
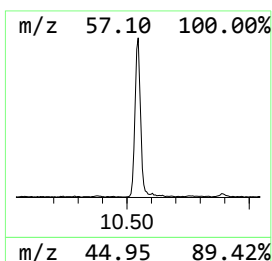
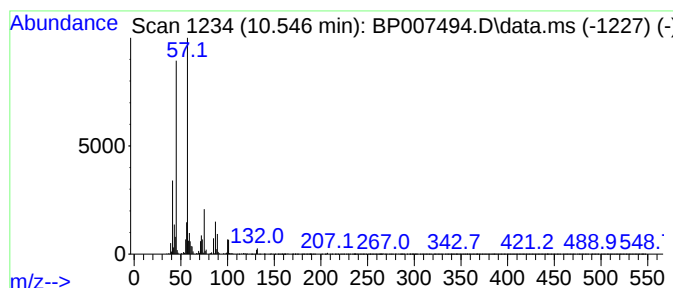
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.546	4.60 ng/ul	345583	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			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	72
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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ALS Vial : 43 Sample Multiplier: 1

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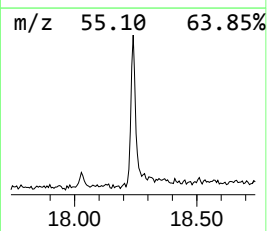
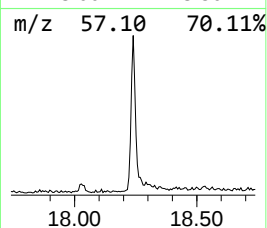
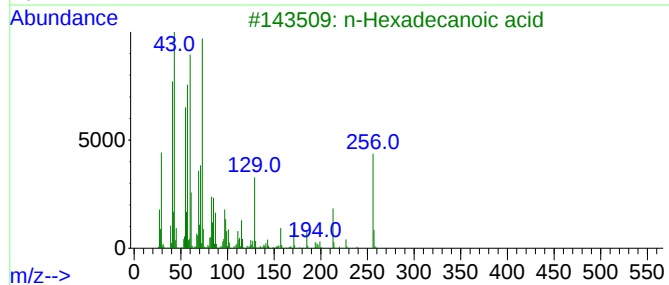
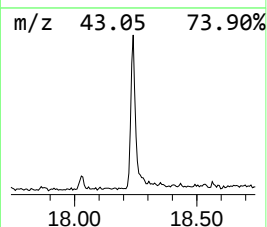
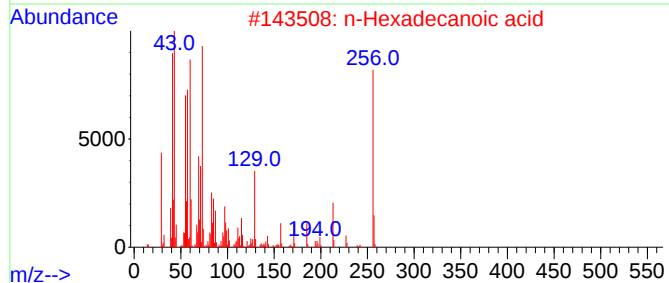
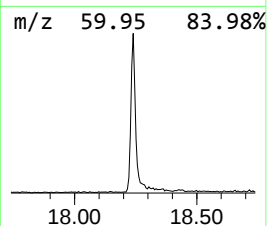
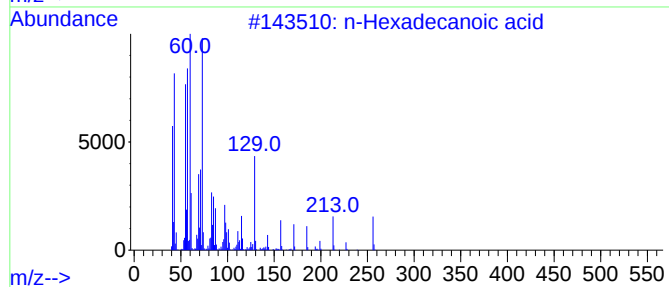
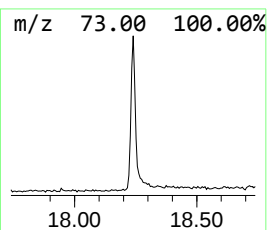
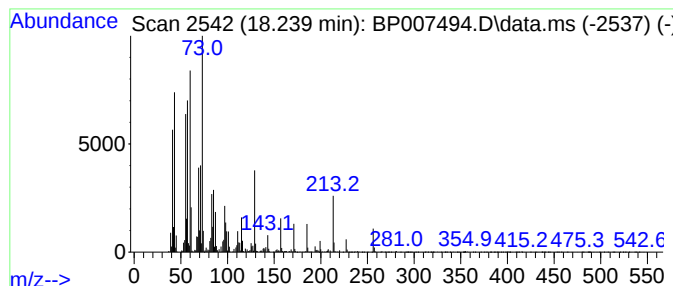
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	3.47 ng/ul	430459	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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



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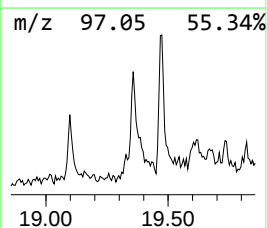
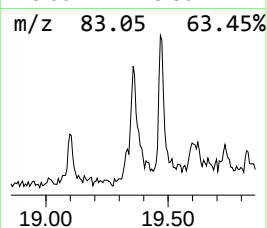
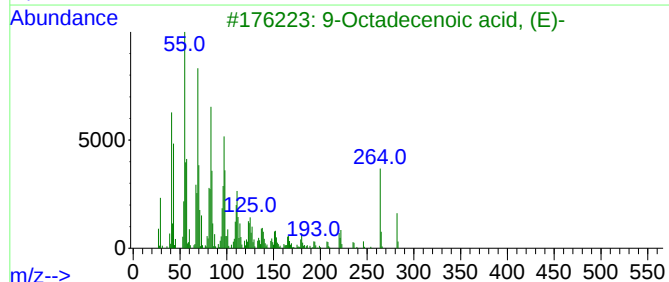
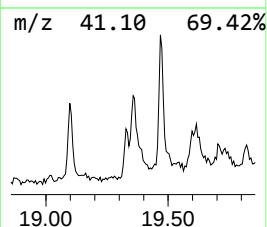
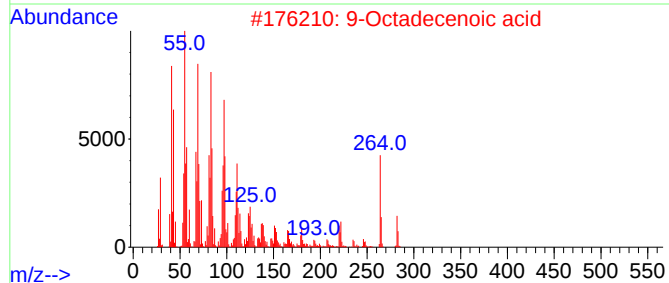
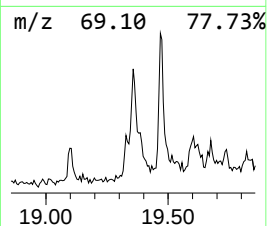
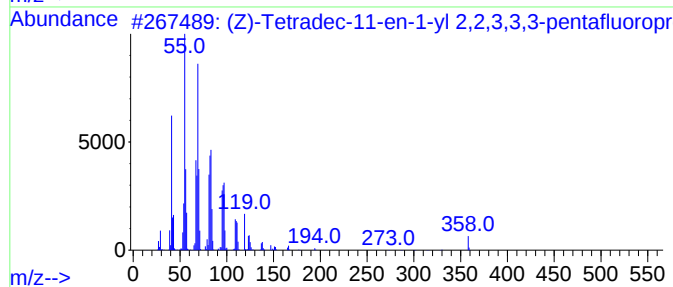
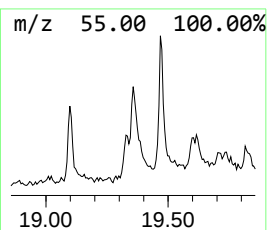
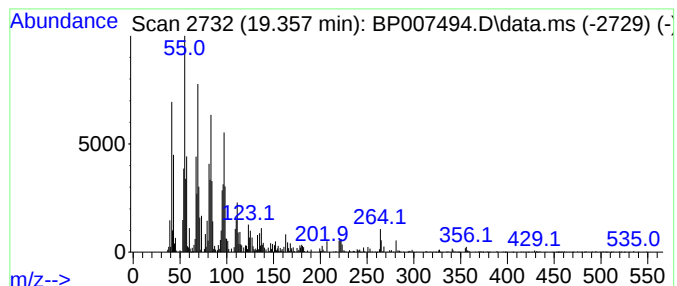
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (Z)-Tetradec-11-en-1-yl 2,2... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.357	2.47 ng/ul	306476	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-Tetradec-11-en-1-yl 2,2,3,3,...	358	C17H27F5O2	1000385-85-5	91		
2	9-Octadecenoic acid	282	C18H34O2	002027-47-6	91		
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	90		
4	Oleyl alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	87		
5	Oleic Acid	282	C18H34O2	000112-80-1	70		



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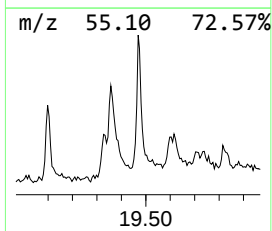
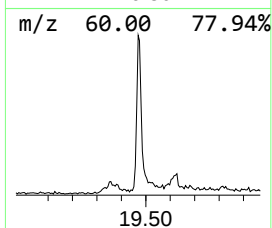
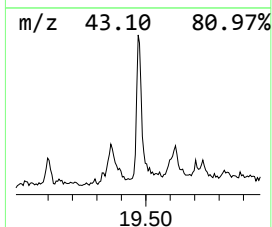
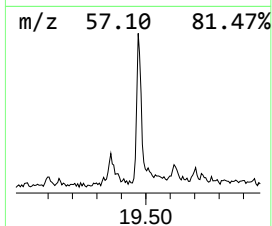
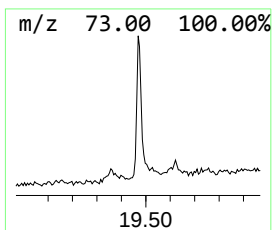
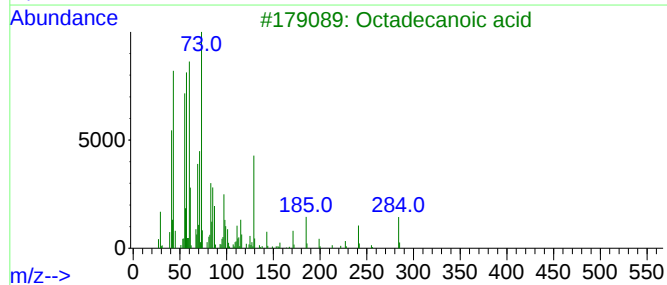
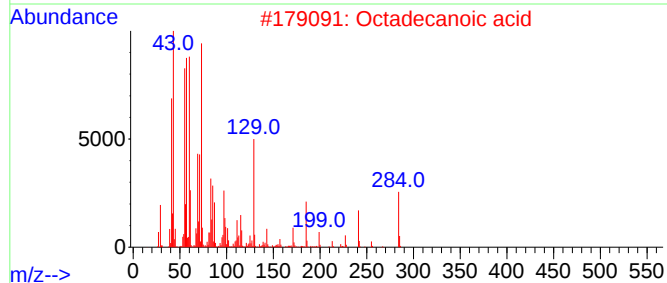
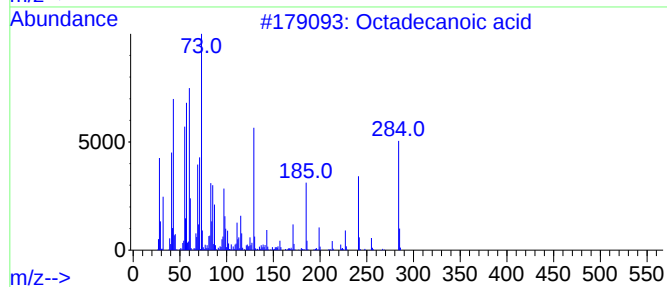
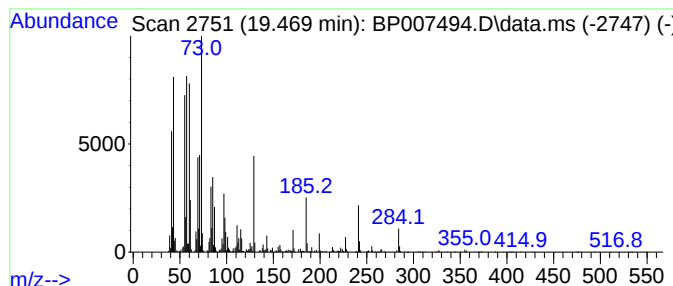
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.469	2.64 ng/ul	386816	Chrysene-d12	21.433

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007494.D
Acq On : 19 Oct 2021 15:21
Operator : CG/JU
Sample : M4196-05
Misc :
ALS Vial : 43 Sample Multiplier: 1

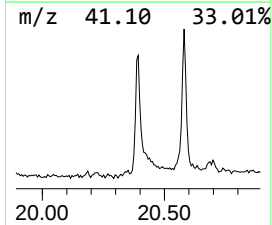
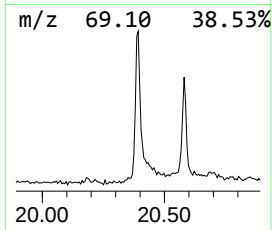
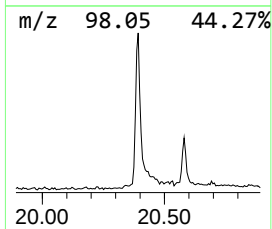
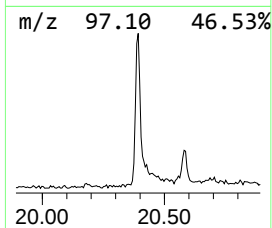
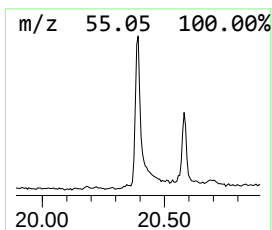
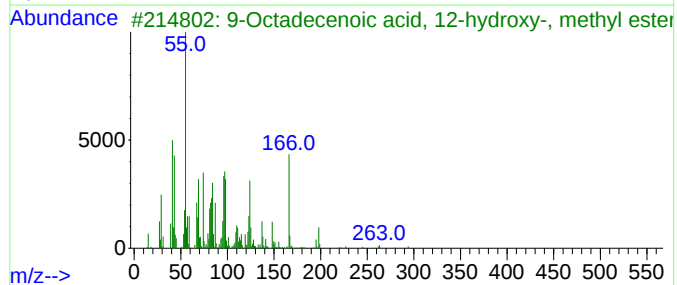
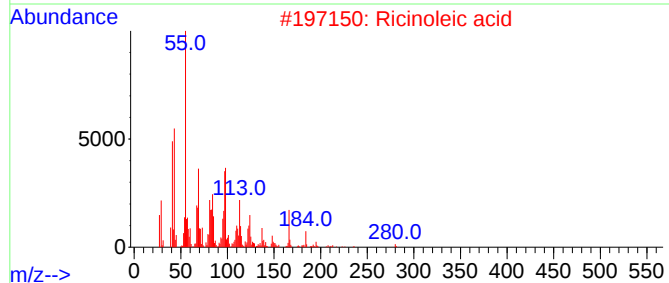
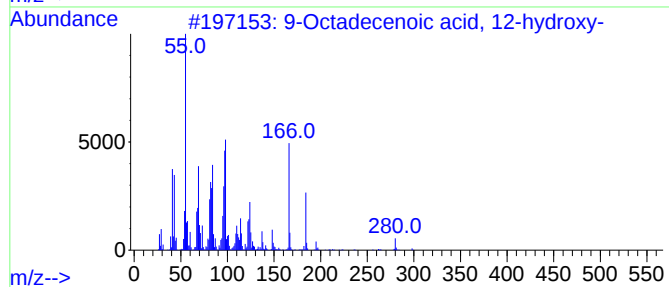
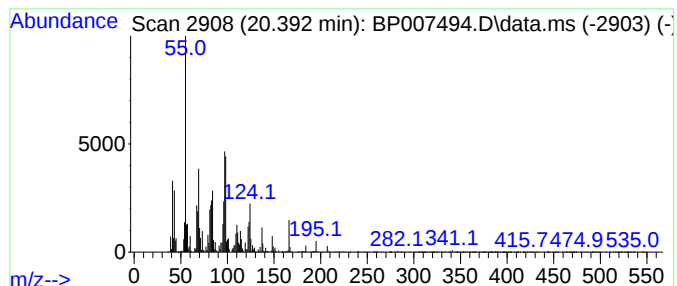
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9-Octadecenoic acid, 12-hyd... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.392	7.04 ng/ul	1031910	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, 12-hydroxy-	298	C18H34O3	007431-95-0	62
2	Ricinoleic acid	298	C18H34O3	000141-22-0	58
3	9-Octadecenoic acid, 12-hydroxy-...	312	C19H36O3	000141-24-2	58
4	Ricinoleic acid	298	C18H34O3	000141-22-0	52
5	9-Octadecenoic acid, 12-hydroxy-...	312	C19H36O3	000141-24-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007494.D
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Instrument :
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ClientSampleId :
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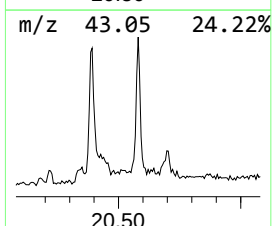
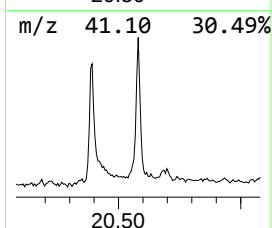
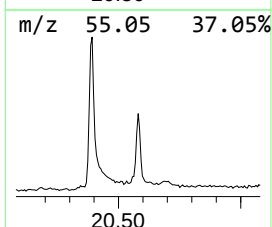
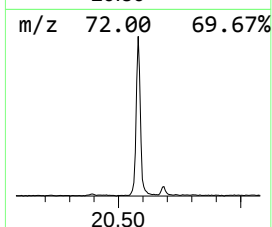
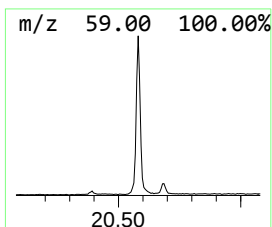
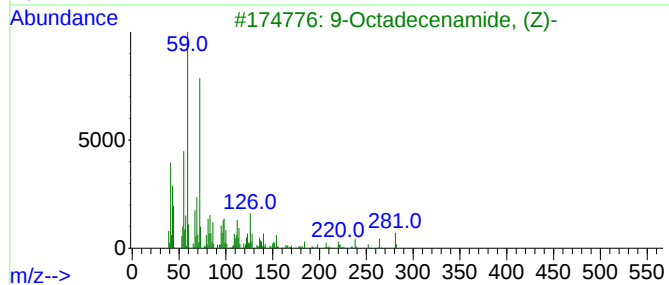
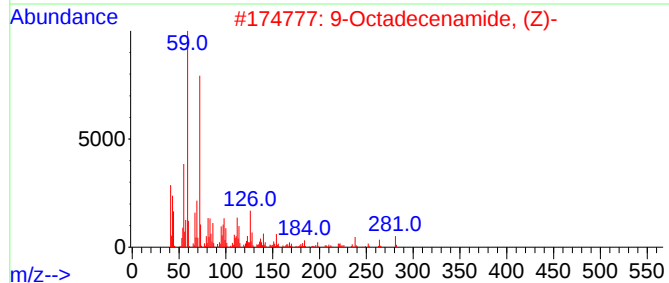
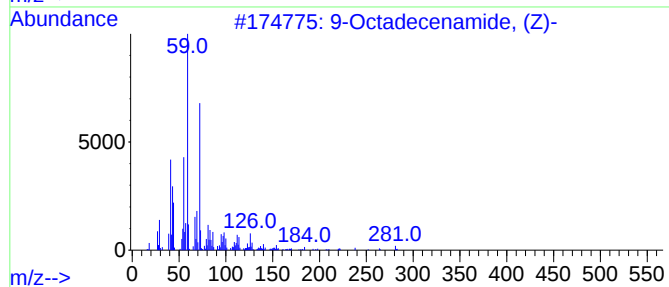
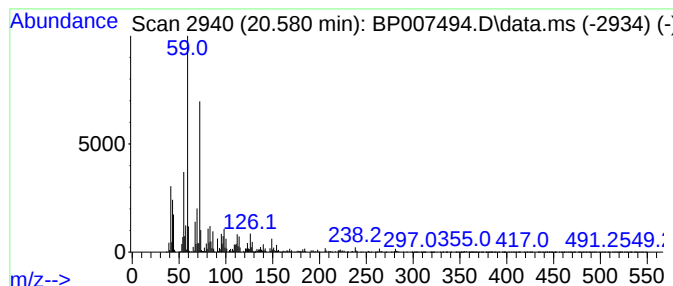
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.580	4.98 ng/ul	730508	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	99
2	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	97
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	97
4	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	95
5	Hexadecanamide			255	C16H33NO	000629-54-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007494.D
Acq On : 19 Oct 2021 15:21
Operator : CG/JU
Sample : M4196-05
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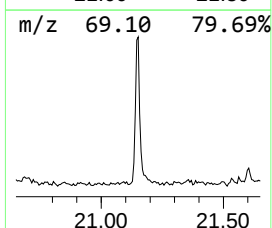
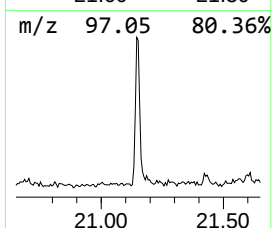
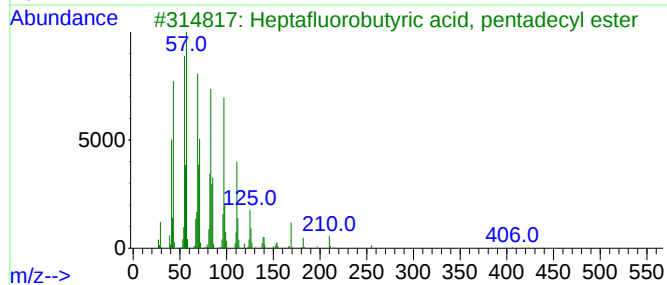
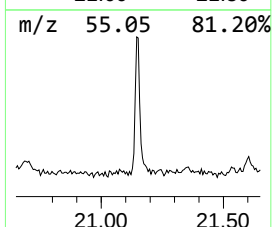
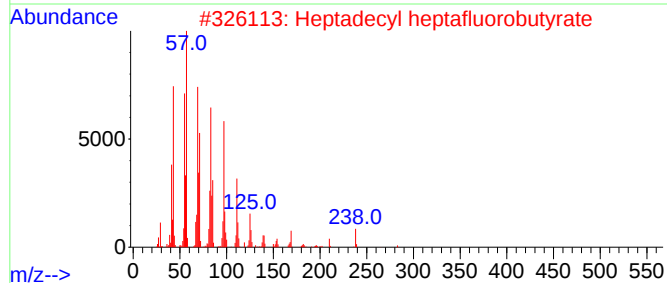
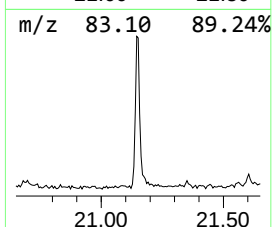
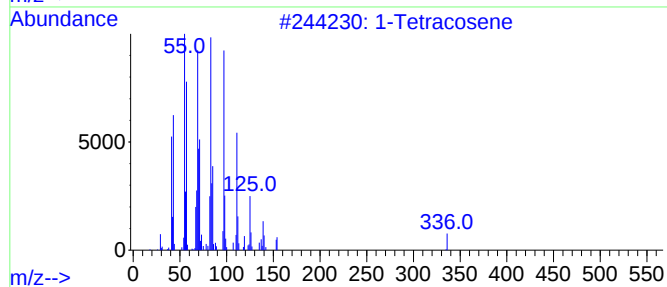
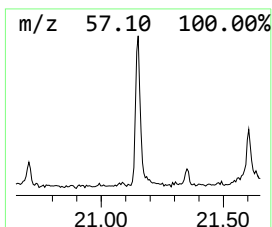
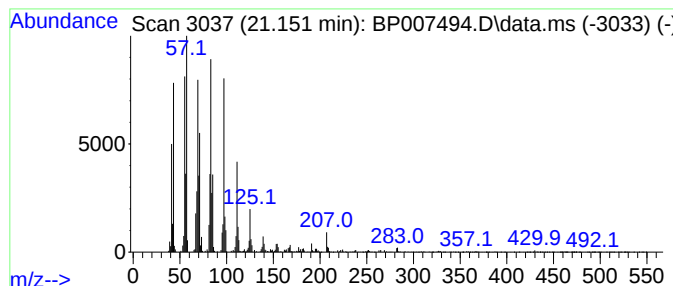
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.151	3.39 ng/ul	497105	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene			336	C24H48	010192-32-2	94
2	Heptadecyl heptafluorobutyrate			452	C21H35F7O2	959085-66-6	94
3	Heptafluorobutyric acid, pentade...			424	C19H31F7O2	959261-23-5	94
4	Nonacos-1-ene			406	C29H58	018835-35-3	94
5	1-Hexacosene			364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007494.D
Acq On : 19 Oct 2021 15:21
Operator : CG/JU
Sample : M4196-05
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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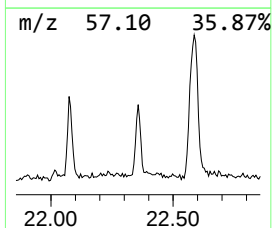
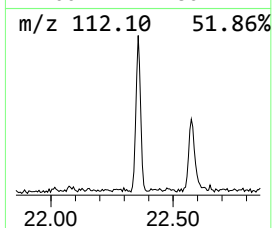
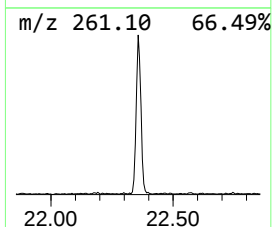
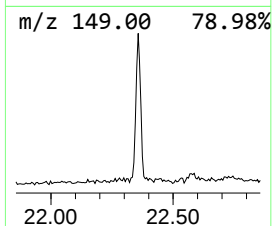
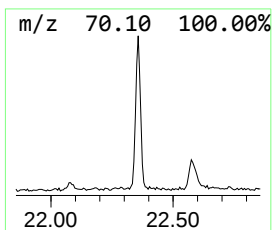
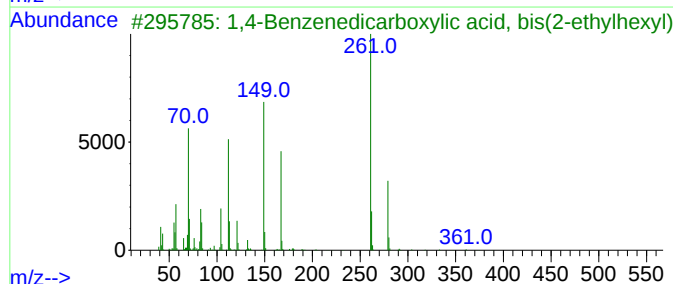
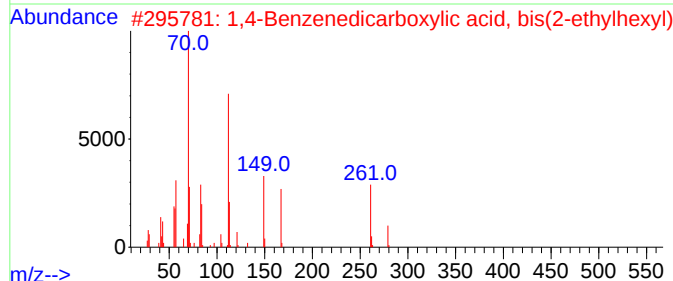
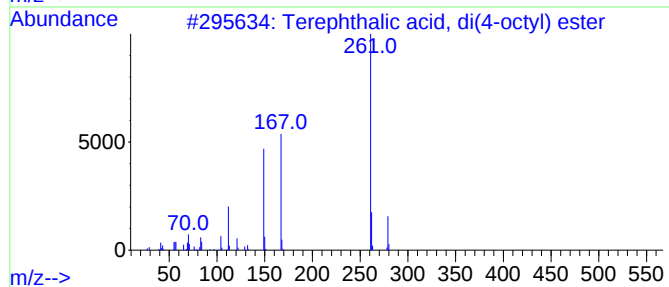
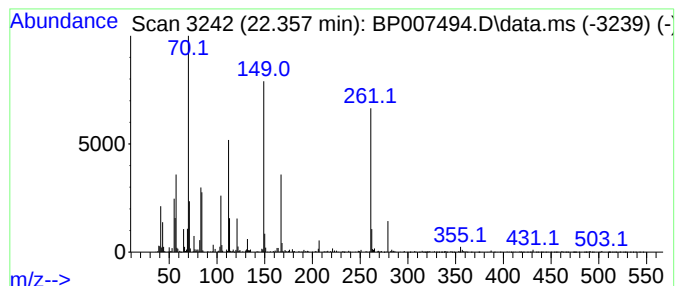
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Terephthalic acid, di(4-oct... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.357	2.09 ng/ul	306761	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Terephthalic acid, di(4-octyl) e...	390	C24H38O4	1000323-74-2	58
2			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	52
3			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	50
4			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	49
5			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007494.D
Acq On : 19 Oct 2021 15:21
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Sample : M4196-05
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Instrument :
BNA_P
ClientSampleId :
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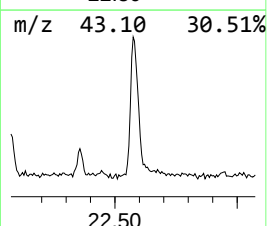
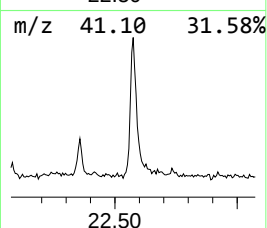
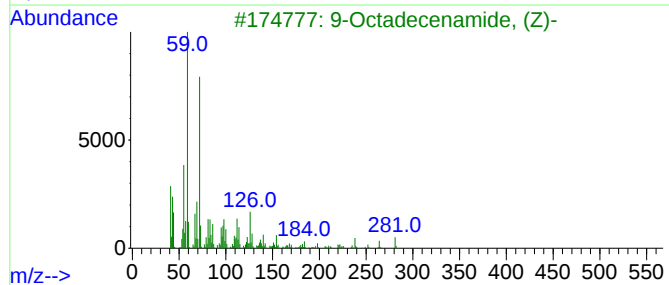
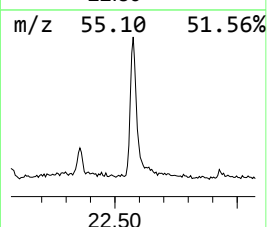
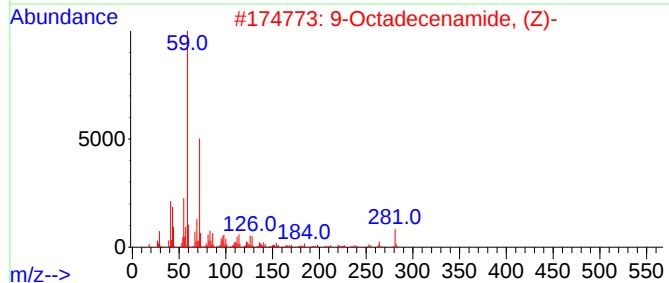
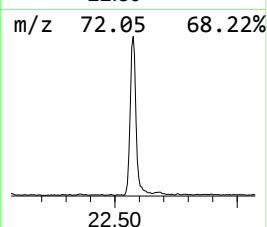
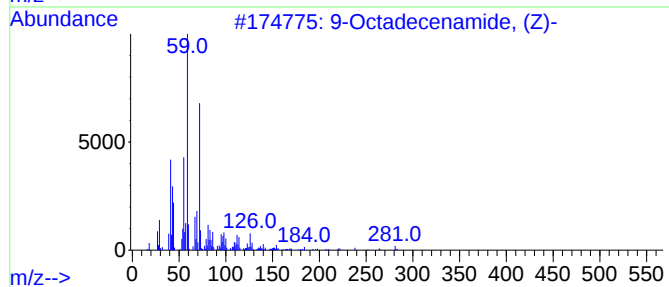
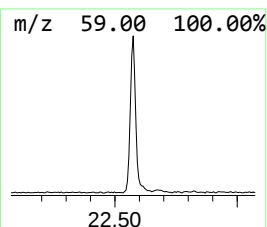
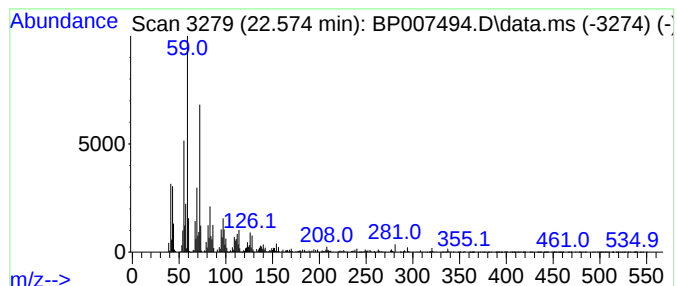
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.574	8.44 ng/ul	1236790	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	93
2	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	93
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	91
4	13-Docosenamide, (Z)-			337	C22H43NO	000112-84-5	90
5	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007494.D
 Acq On : 19 Oct 2021 15:21
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 Sample : M4196-05
 Misc :
 ALS Vial : 43 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
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n-Hexadecanoic ...	18.239	3.5	ng/ul	430459	4	17.345	2483820	20.0
(Z)-Tetradec-11...	19.357	2.5	ng/ul	306476	4	17.345	2483820	20.0
Octadecanoic acid	19.469	2.6	ng/ul	386816	5	21.433	2931760	20.0
9-Octadecenoic ...	20.392	7.0	ng/ul	1031910	5	21.433	2931760	20.0
9-Octadecenamid...	20.580	5.0	ng/ul	730508	5	21.433	2931760	20.0
(DEL) Alkane: S...	21.151	3.4	ng/ul	497105	5	21.433	2931760	20.0
Terephthalic ac...	22.357	2.1	ng/ul	306761	5	21.433	2931760	20.0
13-Docosenamide...	22.574	8.4	ng/ul	1236790	5	21.433	2931760	20.0