LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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: BP007493.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.381	12	16	23	rVB	55145	73480	1.09%	0.103%
2	3.652	59	62	73	rVB	22861	34003	0.51%	0.048%
3	3.793	83	86	108	rVB	588955	875271	13.01%	1.231%
4	4.040	125	128	133	rVB2	13949	21495	0.32%	0.030%
5	4.134	140	144	151	rVB2	13667	18697	0.28%	0.026%
6	4.228	156	160	164	rVB6	6611	8975	0.13%	0.013%
7	5.058	297	301	307	rVB3	9737	15416	0.23%	0.022%
8	6.940	618	621	627	rVB5	9570	14412	0.21%	0.020%
9	7.116	643	651	661	rBV	1086578	1703571	25.33%	2.395%
10	7.281	672	679	689	rVB	811234	1276098	18.97%	1.794%
11	7.487	706	714	721	rBV	1045912	1676531	24.93%	2.357%
12	7.558	723	726	733	rVB	14362	20558	0.31%	0.029%
13	7.958	787	794	801	rBV	713487	1155566	17.18%	1.625%
14	8.022	801	805	812	rVB6	6716	10565	0.16%	0.015%
15	8.663	903	914	922	rBV	1341948	2212366	32.90%	3.111%
16	9.110	982	990	1001	rBV	978924	1615155	24.02%	2.271%
17	9.234	1007	1011	1015	rVB7	4051	6974	0.10%	0.010%
18	9.287	1015	1020	1026	rBV4	18887	30009	0.45%	0.042%
19	9.587	1067	1071	1075	rVB6	6674	9438	0.14%	0.013%
20	9.834	1104	1113	1122	rBV	767685	1310159	19.48%	1.842%
21	10.069	1151	1153	1160	rVB8	6088	11633	0.17%	0.016%
22	10.375	1198	1205	1214	rBV	1344615	2222144	33.04%	3.125%
23	10.540	1227	1233	1249	rBV	425704	673155	10.01%	0.947%
24	10.763	1261	1271	1278	rVB	981151	1633764	24.29%	2.297%
25	10.893	1284	1293	1304	rBV	1245813	2161400	32.14%	3.039%
26	11.557	1401	1406	1410	rVB6	9805	14569	0.22%	0.020%
27	12.251	1521	1524	1529	rVB6	5151	7095	0.11%	0.010%
28	12.346	1535	1540	1547	rVB3	23485	34291	0.51%	0.048%
29	12.704	1594	1601	1608	rBV3	4649	12047	0.18%	0.017%
30	12.963	1638	1645	1650	rBV6	7848	13250	0.20%	0.019%
31	13.210	1682	1687	1693	rVB5	11509	20417	0.30%	0.029%
32	13.440	1720	1726	1730	rBV8	5402	9560	0.14%	0.013%
33	13.504	1733	1737	1741	rVB4	11708	17135	0.25%	0.024%
34	13.563	1743	1747	1752	rVB5	8611	12097	0.18%	0.017%
35	13.934	1805	1810	1815	rBV7	21496	43425	0.65%	0.061%
36	13.998	1815	1821	1835	rVB	2165865	3092349	45.98%	4.348%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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	14.175	1847	1851	1857	rVB6	8484	12745	0.19%	0.018%
38	14.281	1859	1869	1877	rBV	2450971	3871615	57.57%	5.444%
39	14.363	1879	1883	1887	rVB6	10597	15411	0.23%	0.022%
40	14.592	1913	1922	1931	rBV	1519980	2213247	32.91%	3.112%
41	14.775	1944	1953	1965	rBV	1112402	1647883	24.50%	2.317%
42	15.004	1988	1992	1998	rBV7	22293	35893	0.53%	0.050%
43	15.257	2028	2035	2040	rBV7	7550	19087	0.28%	0.027%
44	15.410	2056	2061	2066	rBV3	10237	16483	0.25%	0.023%
45	15.587	2084	2091	2098	rBV	3252521	4727206	70.29%	6.647%
46	15.692	2103	2109	2121	rBV	687373	976486	14.52%	1.373%
47	15.828	2126	2132	2136	rVV2	39011	59040	0.88%	0.083%
48	15.951	2151	2153	2159	rVB7	4797	7560	0.11%	0.011%
49	16.034	2163	2167	2171	rVB7	8894	14381	0.21%	0.020%
50	16.151	2182	2187	2194	rBV7	6951	14353	0.21%	0.020%
51	16.257	2202	2205	2212	rVB8	6460	13354	0.20%	0.019%
52	16.386	2220	2227	2230	rVB9	13010	24161	0.36%	0.034%
53	16.592	2260	2262	2268	rBV7	5323	9989	0.15%	0.014%
54	16.751	2284	2289	2294	rVB6	19151	28030	0.42%	0.039%
55	16.957	2318	2324	2325	rBV6	4506	8604	0.13%	0.012%
56	17.022	2332	2335	2341	rVB7	6326	10609	0.16%	0.015%
57	17.128	2351	2353	2357	rVB2	9073	11839	0.18%	0.017%
58	17.345	2383	2390	2397	rBV	1916672	2731200	40.61%	3.840%
59	17.445	2399	2407	2416	rBV	3556241	5180571	77.03%	7.285%
60	17.533	2417	2422	2427	rVV3	43837	65188	0.97%	0.092%
61	17.639	2438	2440	2449	rVB9	5945	12421	0.18%	0.017%
62	17.710	2449	2452	2455	rBV2	7004	9696	0.14%	0.014%
63	18.028	2502	2506	2512	rVB	40436	55658	0.83%	0.078%
64	18.110	2518	2520	2525	rVB6	9524	11006	0.16%	0.015%
65	18.239	2537	2542	2550	rBV	462316	635767	9.45%	0.894%
66	19.010	2671	2673	2677	rVB4	14240	15898	0.24%	0.022%
67	19.098	2680	2688	2696	rBV	300140	452100	6.72%	0.636%
68	19.251	2710	2714	2718	rBV2	51643	70542	1.05%	0.099%
69	19.327	2722	2727	2729	rVV2	118216	169173	2.52%	0.238%
70	19.357	2729	2732	2743	rVV2	179405	379437	5.64%	0.534%
71	19.469	2747	2751	2763	rVV	268127	407660	6.06%	0.573%
72	19.675	2780	2786	2791	rBV2	4530497	6293633	93.58%	8.850%
73	19.822	2808	2811	2818	rVB5	37961	52462	0.78%	0.074%
74	20.392	2903	2908	2925	rBV	763912	1394571	20.74%	1.961%
75	20.580	2934	2940	2948	rBV	1064156	1301281	19.35%	1.830%
76	21.145	3032	3036	3047	rBV	584317	844736	12.56%	1.188%

LSC Area Percent Report

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Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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

77	21.427	3079	3084	3091	rBV	2419604	3343164	49.71%	4.701%
78	22.016	3181	3184	3190	rVB5	115100	152598	2.27%	0.215%
79	22.574	3274	3279	3287	rBV	901035	1476897	21.96%	2.077%
80	23.727	3467	3475	3483	rBV2	3292885	6725465	100.00%	9.457%
81	23.886	3495	3502	3511	rVB2	1616904	3531045	52.50%	4.965%

Sum of corrected areas: 71117215

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
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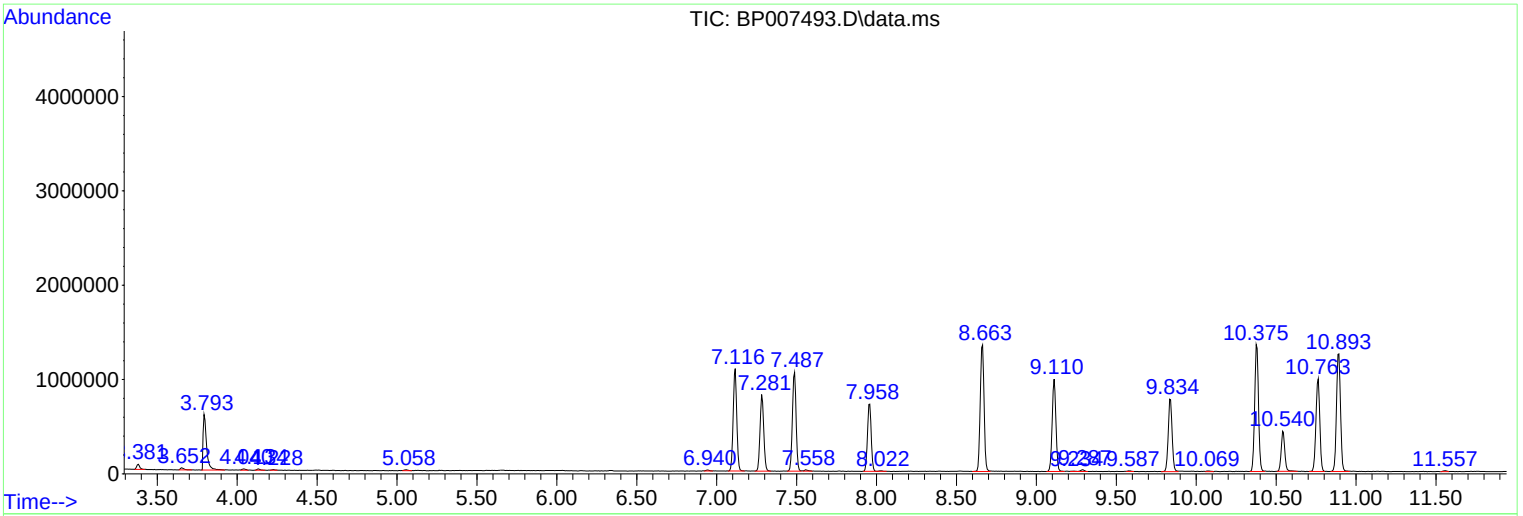
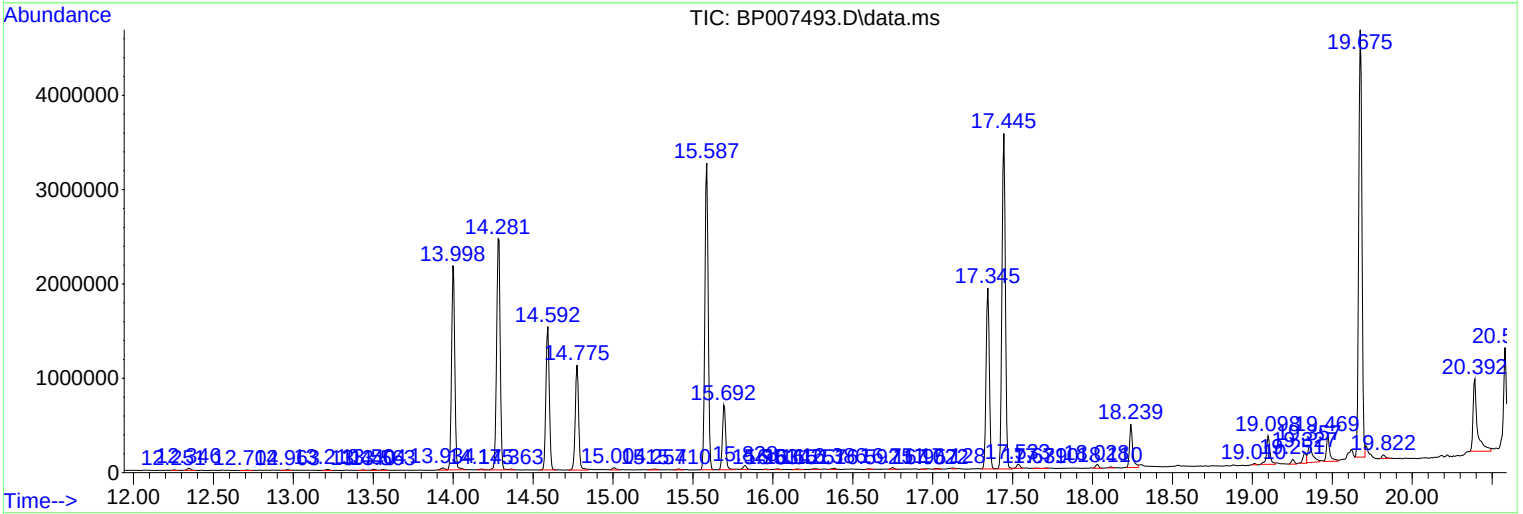
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

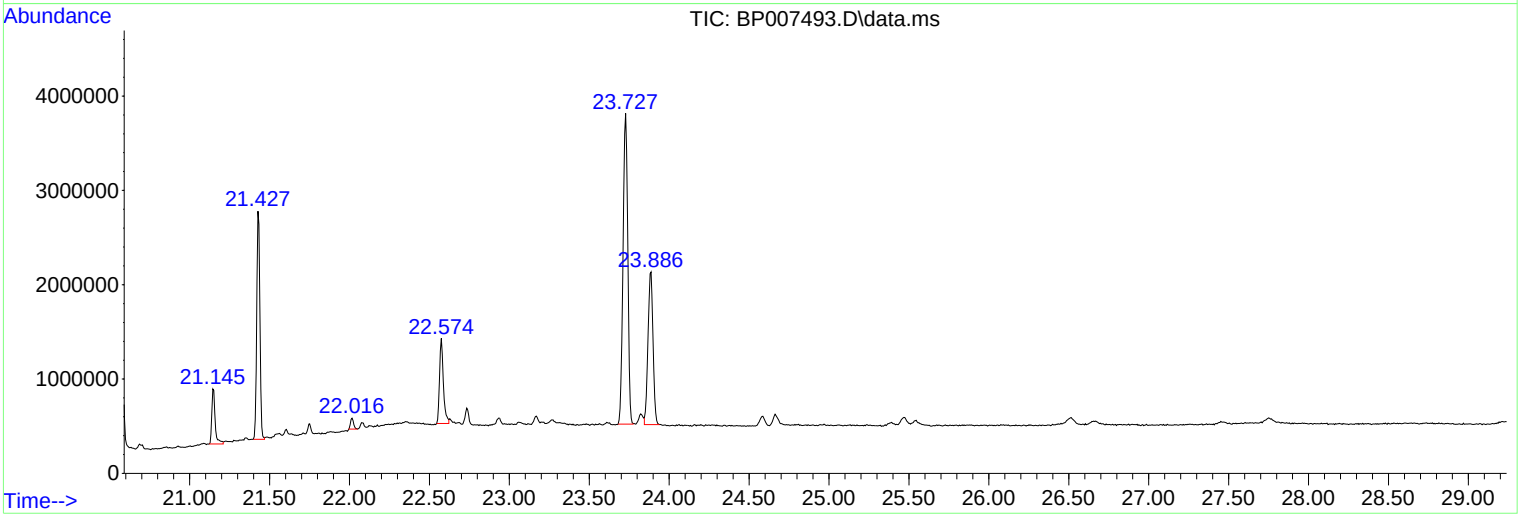
Instrument :
BNA_P
ClientSampleId :
JDGA8

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 : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

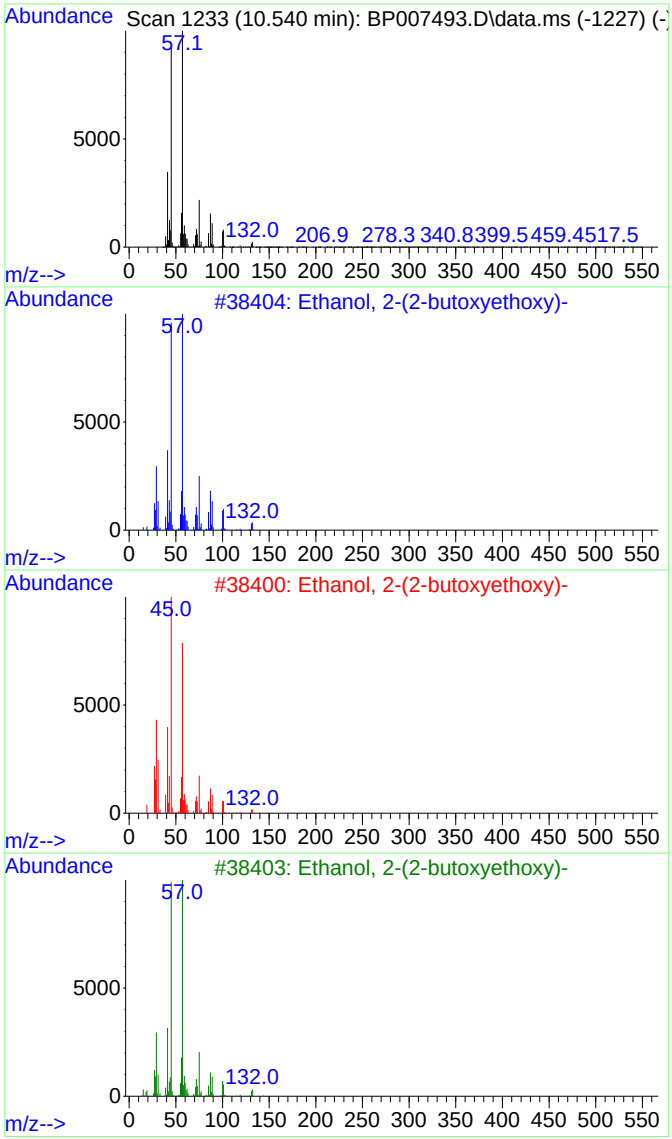
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

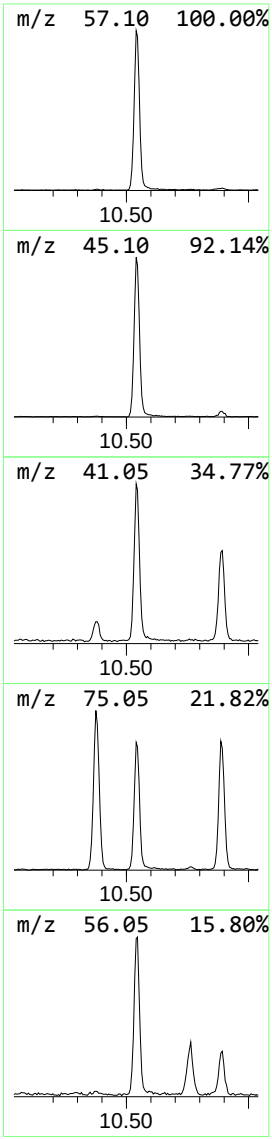
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	8.24 ng/ul	673155	Naphthalene-d8	10.763

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	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





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Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA8

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

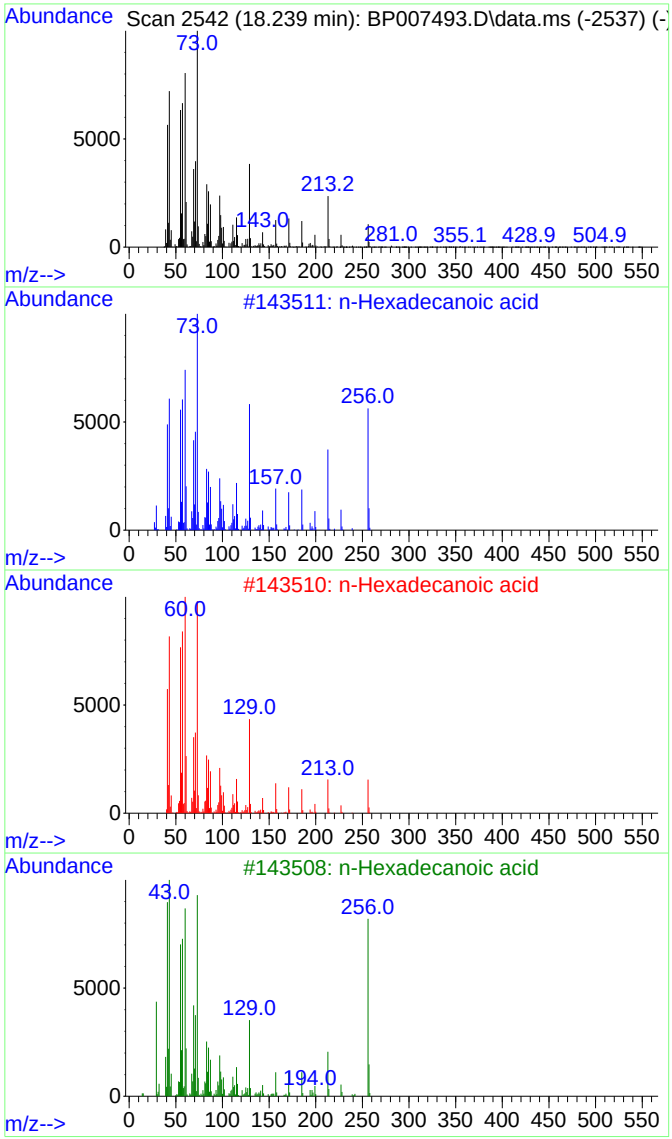
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

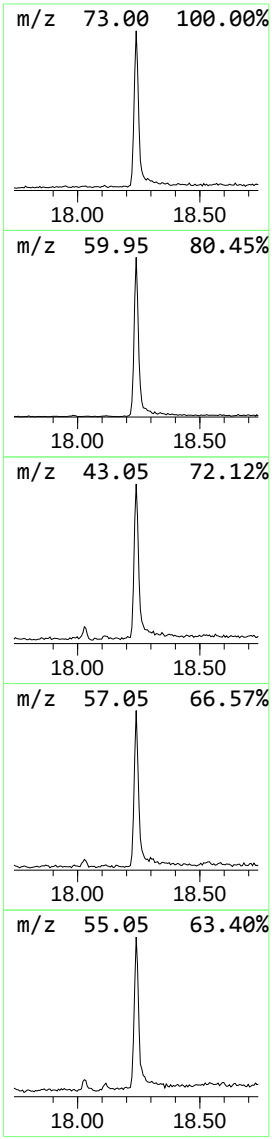
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	4.66 ng/ul	635767	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	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

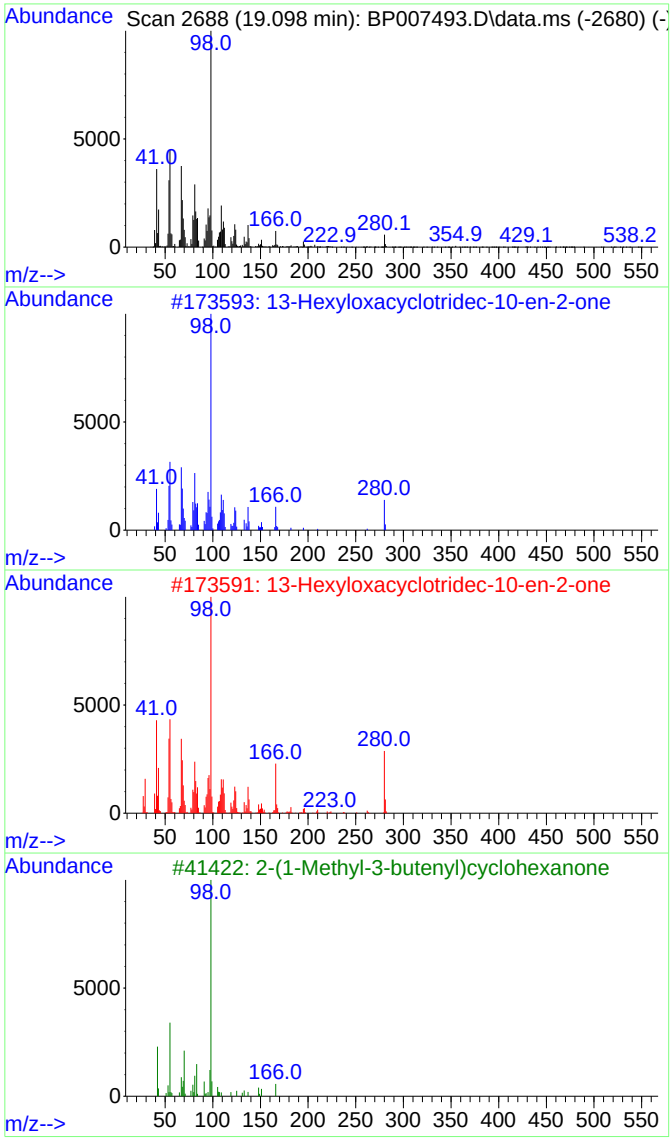
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

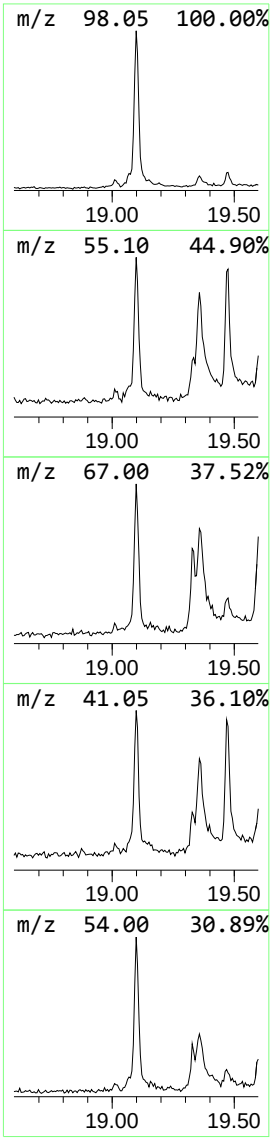
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 13-Hexyloxacyclotridec-10-e... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.098	3.31 ng/ul	452100	Phenanthrene-d10	17.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Hexyloxacyclotridec-10-en-2-one	280	C18H32O2	127062-51-5	99
2	13-Hexyloxacyclotridec-10-en-2-one	280	C18H32O2	127062-51-5	95
3	2-(1-Methyl-3-butenyl)cyclohexanone	166	C11H18O	1000424-67-7	43
4	(Cyclopentylamino)acetic acid, m...	157	C8H15NO2	190904-15-5	43
5	4-Chloro-N-[2-(1-piperidinyl)eth...	266	C14H19ClN2O	189203-75-6	43





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Acq On : 19 Oct 2021 14:47
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ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA8

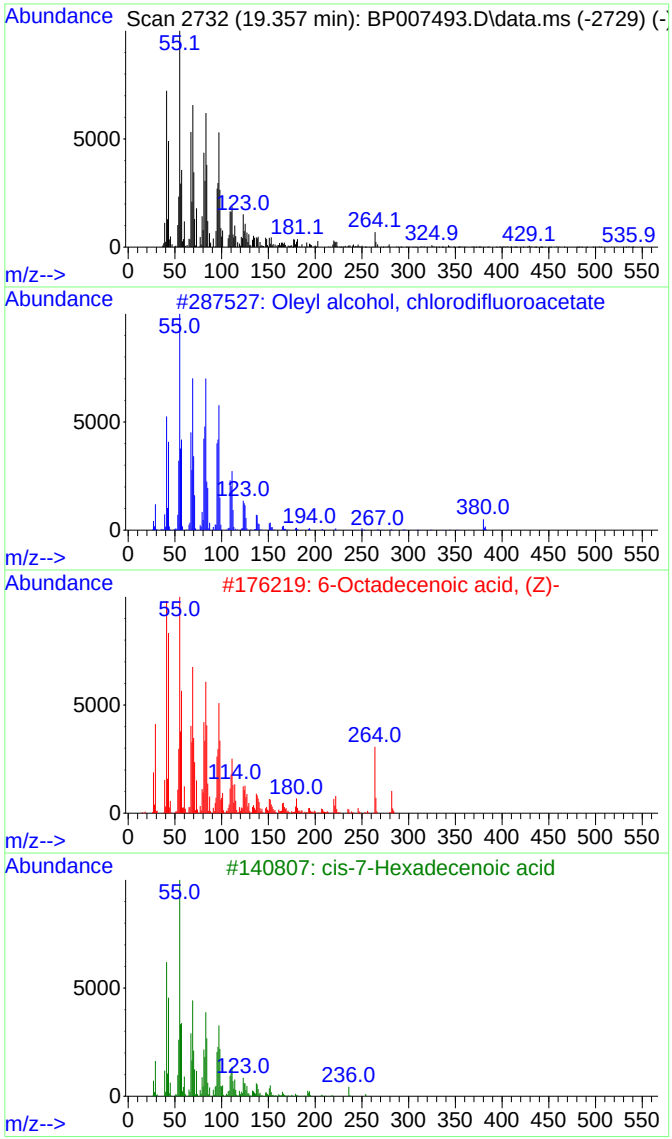
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

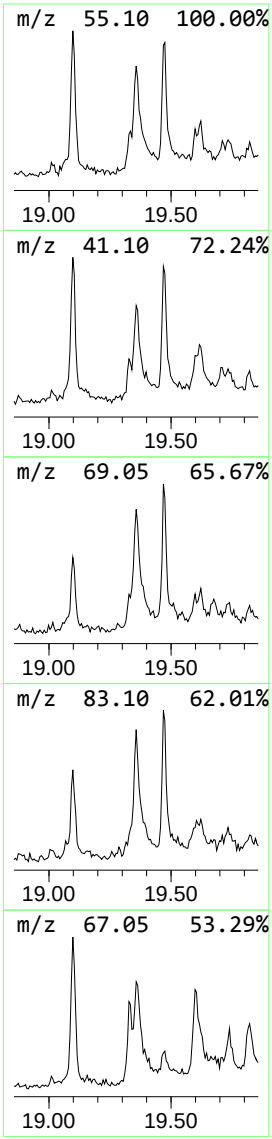
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Oleyl alcohol, chlorodifluo... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.357	2.78 ng/ul	379437	Phenanthrene-d10	17.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleyl alcohol, chlorodifluoroace...	380	C20H35ClF2O2	1000447-47-7	96
2	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	87
3	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	83
4	9-Octadecenoic acid	282	C18H34O2	002027-47-6	81
5	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	81





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Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

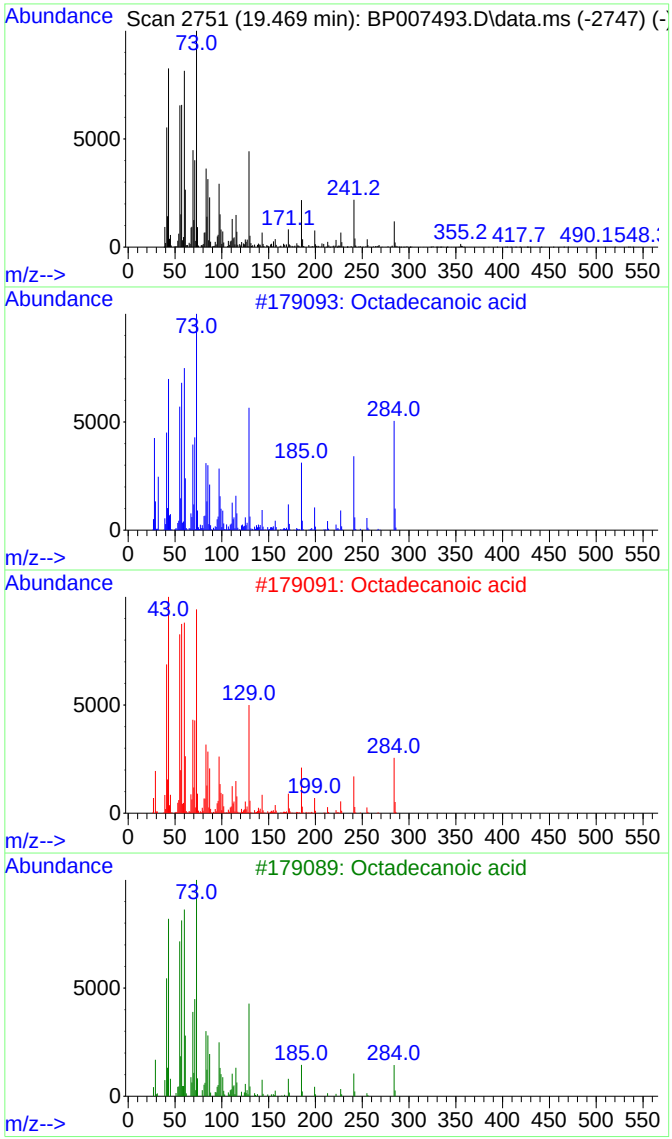
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

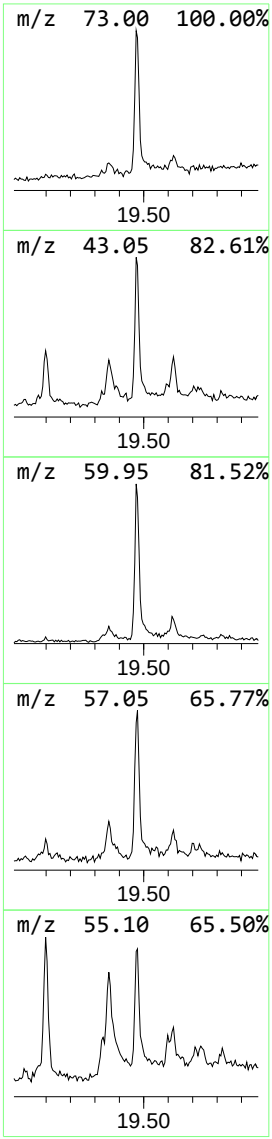
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.469	2.44 ng/ul	407660	Chrysene-d12	21.427

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	96
5	Octadecanoic acid	284	C18H36O2	000057-11-4	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA8

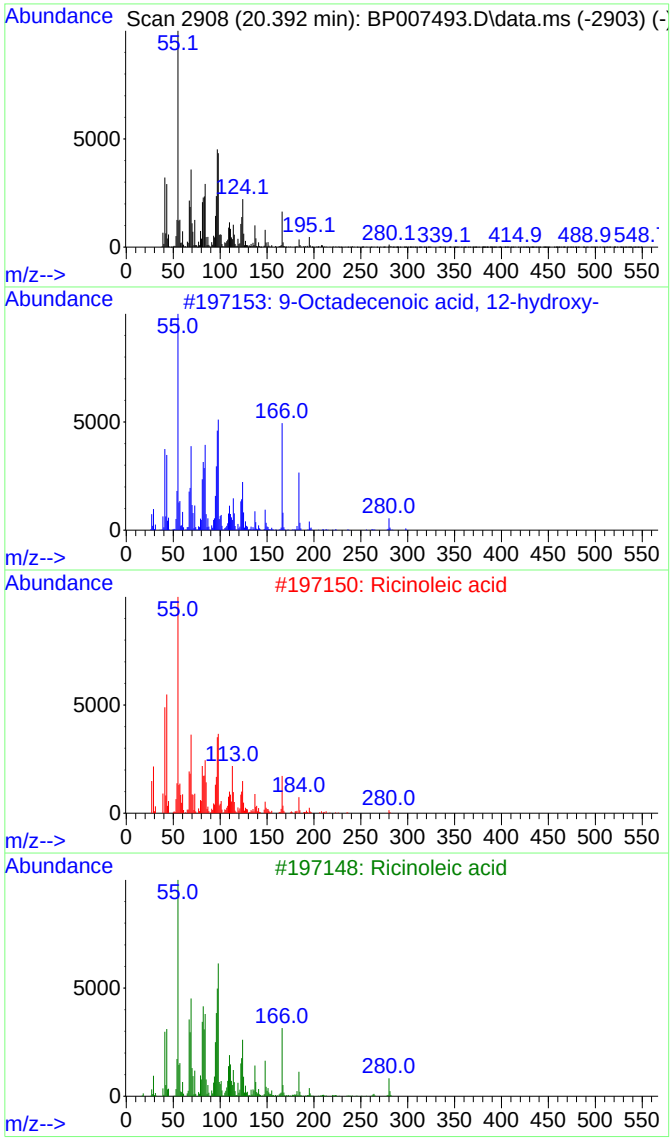
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

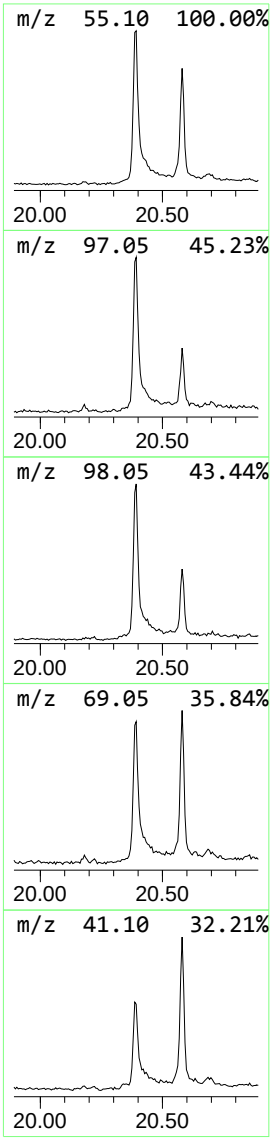
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 9-Octadecenoic acid, 12-hyd... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.392	8.34 ng/ul	1394570	Chrysene-d12	21.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, 12-hydroxy-	298	C18H34O3	007431-95-0	87
2	Ricinoleic acid	298	C18H34O3	000141-22-0	87
3	Ricinoleic acid	298	C18H34O3	000141-22-0	68
4	Ricinoleic acid	298	C18H34O3	000141-22-0	64
5	Chloromethyl 9-chloroundecanoate	268	C12H22Cl2O2	080418-95-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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

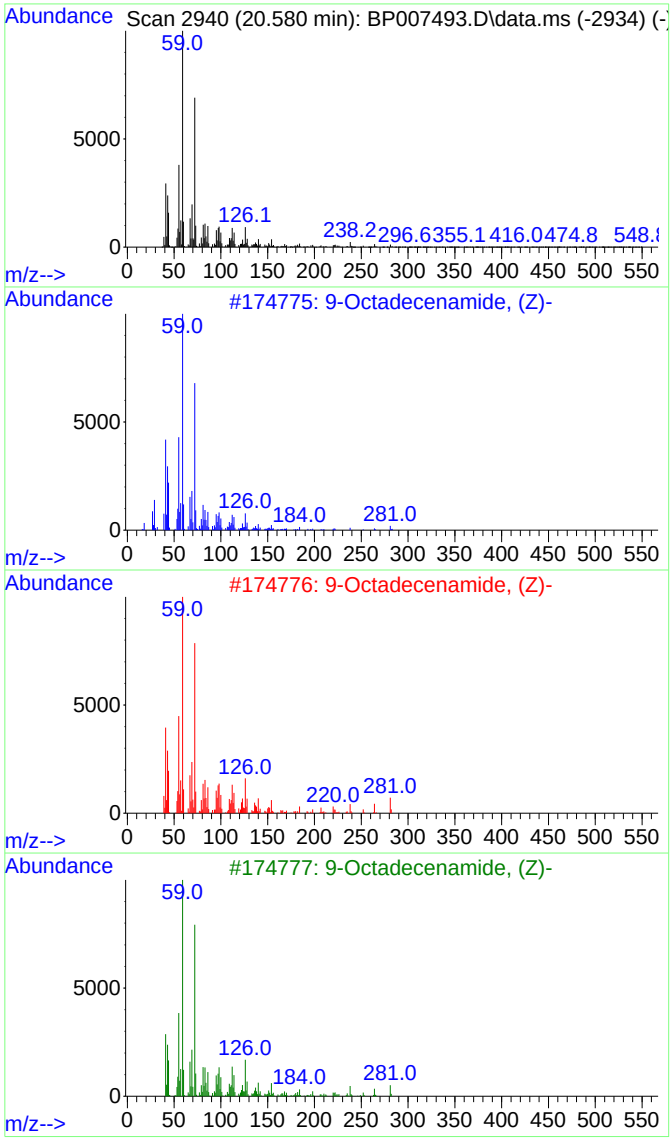
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

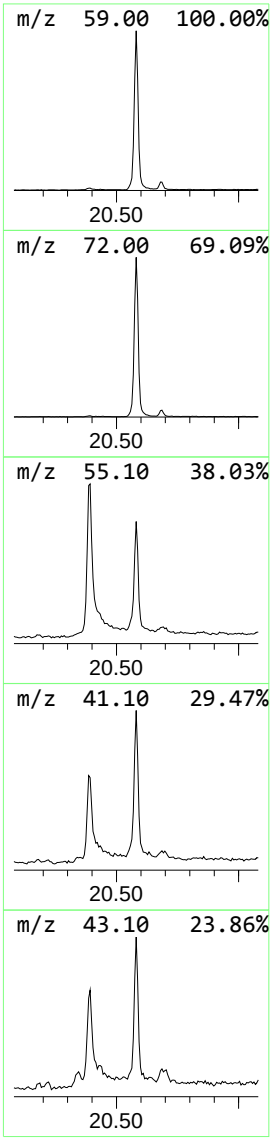
****

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.580	7.78 ng/ul	1301280	Chrysene-d12	21.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	96
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	95
4	Palmitoleamide	253	C16H31NO	106010-22-4	72
5	Dodecanamide	199	C12H25NO	001120-16-7	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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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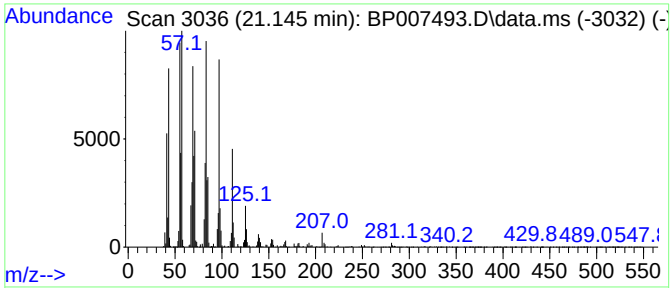
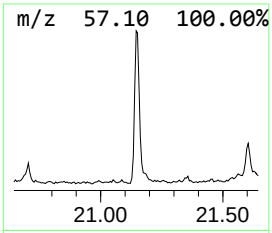
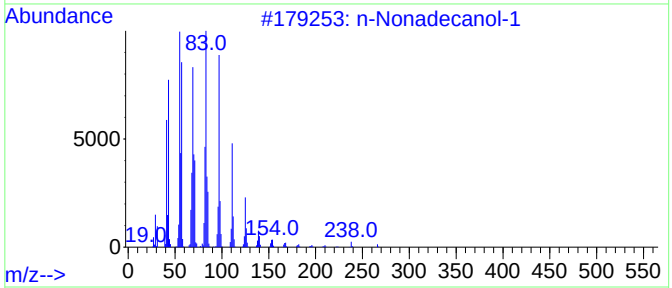
Peak Number 8 n-Nonadecanol-1 Concentration Rank 5

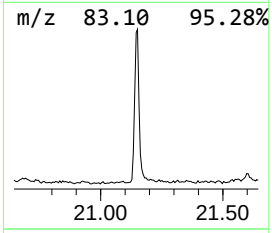
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	5.05 ng/ul	844736	Chrysene-d12	21.427

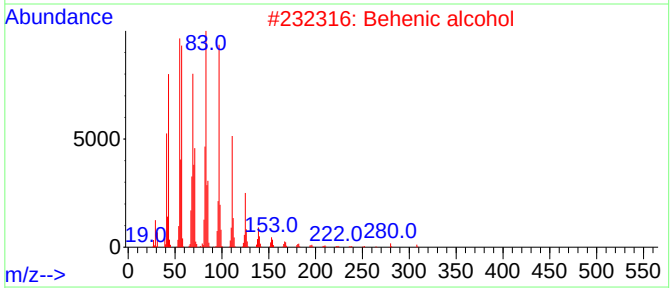
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4	Pentacos-1-ene	350	C25H50	016980-85-1	94
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94

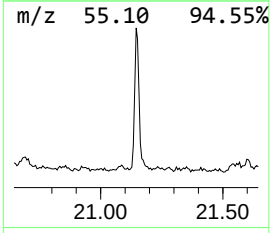


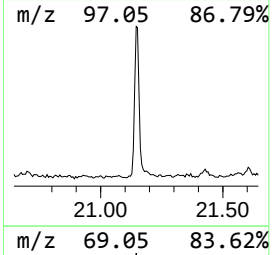


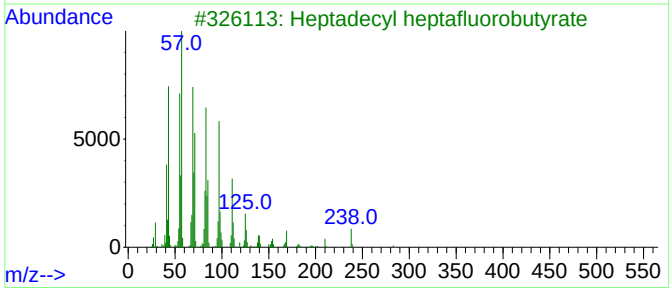


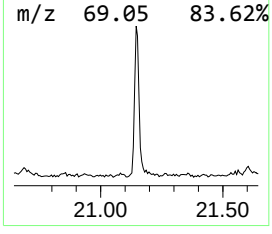












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

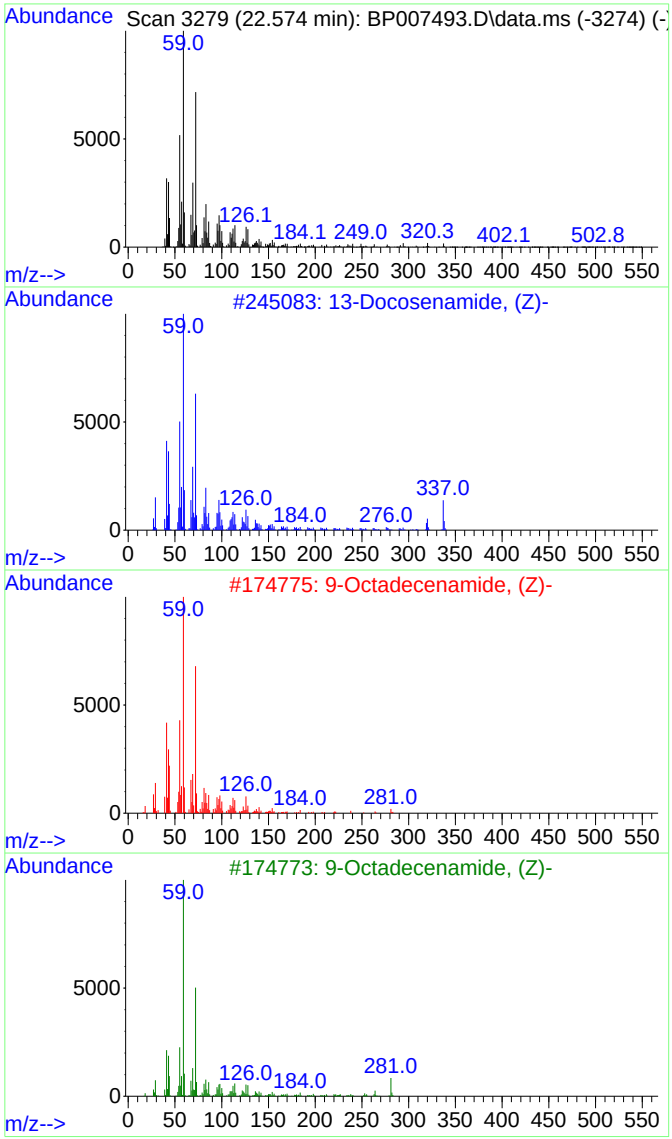
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

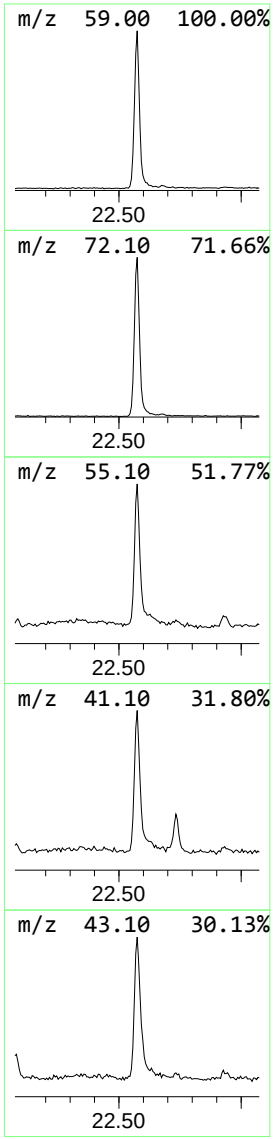
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.574	8.84 ng/ul	1476900	Chrysene-d12	21.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
4	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	76
5	Tetradecanamide	227	C14H29NO	000638-58-4	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007493.D
Acq On : 19 Oct 2021 14:47
Operator : CG/JU
Sample : M4196-07
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGA8

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Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.540	8.2	ng/ul	673155	2	10.763	1633760	20.0
n-Hexadecanoic ...	18.239	4.7	ng/ul	635767	4	17.345	2731200	20.0
13-Hexyloxacycl...	19.098	3.3	ng/ul	452100	4	17.345	2731200	20.0
Oleyl alcohol, ...	19.357	2.8	ng/ul	379437	4	17.345	2731200	20.0
Octadecanoic acid	19.469	2.4	ng/ul	407660	5	21.427	3343160	20.0
9-Octadecenoic ...	20.392	8.3	ng/ul	1394570	5	21.427	3343160	20.0
9-Octadecenamid...	20.580	7.8	ng/ul	1301280	5	21.427	3343160	20.0
n-Nonadecanol-1	21.145	5.0	ng/ul	844736	5	21.427	3343160	20.0
13-Docosenamide...	22.574	8.8	ng/ul	1476900	5	21.427	3343160	20.0