LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

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-BP120123.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP018738.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	26	30	40	rVB	61837	86649	1.42%	0.124%
2	3.558	77	80	90	rVB4	19008	32628	0.53%	0.047%
3	3.681	97	101	119	rVB	657361	1022157	16.75%	1.468%
4	3.958	131	148	153	rBV	277768	824290	13.51%	1.184%
5	4.011	155	157	163	rVB	19024	25612	0.42%	0.037%
6	4.817	290	294	298	rBV7	4810	7287	0.12%	0.010%
7	4.940	312	315	319	rBV3	4340	8577	0.14%	0.012%
8	5.152	344	351	359	rVB	278743	429223	7.03%	0.616%
9	5.287	369	374	379	rBV2	13308	22564	0.37%	0.032%
10	5.340	379	383	388	rVB2	27226	39336	0.64%	0.056%
11	6.381	556	560	564	rBV6	4375	7844	0.13%	0.011%
12	6.864	632	642	648	rBV	114965	229819	3.77%	0.330%
13	7.011	661	667	680	rVB	847402	1342180	22.00%	1.927%
14	7.175	688	695	706	rVB	695903	1095846	17.96%	1.574%
15	7.369	721	728	738	rBV	848816	1322570	21.68%	1.899%
16	7.452	739	742	749	rVB4	11272	20297	0.33%	0.029%
17	7.728	781	789	796	rVB	305116	471574	7.73%	0.677%
18	7.840	801	808	811	rVV	833218	1335069	21.88%	1.917%
19	7.875	811	814	824	rVB	876720	1293546	21.20%	1.857%
20	8.193	860	868	875	rBV	797068	1238532	20.30%	1.778%
21	8.552	922	929	942	rBV	1052780	1620749	26.56%	2.327%
22	8.999	998	1005	1014	rBV	731866	1182952	19.39%	1.699%
23	9.240	1043	1046	1053	rVB7	8076	13423	0.22%	0.019%
24	9.722	1121	1128	1134	rBV	555701	885173	14.51%	1.271%
25	9.781	1134	1138	1142	rVB4	14616	23582	0.39%	0.034%
26	10.257	1212	1219	1238	rBV	976000	1628950	26.70%	2.339%
27	10.422	1243	1247	1250	rVB6	4530	6601	0.11%	0.009%
28	10.499	1254	1260	1273	rVB	426132	704847	11.55%	1.012%
29	10.634	1273	1283	1296	rBV	1074871	1788959	29.32%	2.569%
30	10.775	1296	1307	1319	rBV	879143	1584685	25.97%	2.275%
31	11.016	1343	1348	1355	rVB	51231	87733	1.44%	0.126%
32	11.446	1415	1421	1423	rBV5	3992	6814	0.11%	0.010%
33	11.810	1480	1483	1486	rBV5	5764	7968	0.13%	0.011%
34	12.063	1522	1526	1530	rVB7	7726	13157	0.22%	0.019%
35	12.222	1548	1553	1558	rBV	15909	26822	0.44%	0.039%
36	12.440	1586	1590	1600	rBV9	9477	17998	0.29%	0.026%

LSC Area Percent Report

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Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

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-BP120123.MA.M
Title : SVOA CALIBRATION

37	12.663	1619	1628	1637	rBV2	37765	78484	1.29%	0.113%
38	13.269	1726	1731	1736	rBV	63180	98518	1.61%	0.141%
39	13.375	1743	1749	1757	rBV4	20279	37862	0.62%	0.054%
40	13.440	1757	1760	1766	rVB8	8204	15242	0.25%	0.022%
41	13.498	1766	1770	1772	rBV5	4902	6301	0.10%	0.009%
42	13.887	1828	1836	1847	rBV	1803965	2479068	40.63%	3.560%
43	14.163	1875	1883	1895	rBV	2067679	2955560	48.44%	4.244%
44	14.469	1929	1935	1944	rVB	1635320	2348327	38.49%	3.372%
45	14.669	1963	1969	1983	rBV	704086	1115911	18.29%	1.602%
46	14.787	1985	1989	1993	rVB6	11768	20457	0.34%	0.029%
47	14.893	2002	2007	2013	rVB9	12315	22014	0.36%	0.032%
48	15.463	2097	2104	2112	rBV	2745763	3974337	65.14%	5.707%
49	15.581	2118	2124	2139	rBV	615493	838286	13.74%	1.204%
50	15.704	2141	2145	2153	rVB5	13631	24466	0.40%	0.035%
51	16.187	2225	2227	2231	rVB5	6682	8262	0.14%	0.012%
52	16.251	2233	2238	2243	rBV9	4610	9417	0.15%	0.014%
53	16.469	2272	2275	2281	rVB5	5987	10484	0.17%	0.015%
54	16.616	2298	2300	2307	rBV8	8011	12877	0.21%	0.018%
55	17.004	2361	2366	2370	rBV5	4912	9362	0.15%	0.013%
56	17.216	2396	2402	2409	rBV	2087566	2977858	48.81%	4.276%
57	17.316	2413	2419	2429	rVV	3215373	4465133	73.18%	6.412%
58	17.416	2433	2436	2441	rVB6	24252	31738	0.52%	0.046%
59	17.863	2508	2512	2521	rVB6	20837	45691	0.75%	0.066%
60	18.110	2549	2554	2559	rBV	304540	407968	6.69%	0.586%
61	18.292	2581	2585	2588	rBV5	33798	45841	0.75%	0.066%
62	18.498	2614	2620	2627	rBV	477152	644397	10.56%	0.925%
63	18.557	2627	2630	2638	rVB6	23051	37041	0.61%	0.053%
64	18.792	2666	2670	2677	rVB3	93203	128804	2.11%	0.185%
65	19.010	2703	2707	2720	rVB	377541	513177	8.41%	0.737%
66	19.133	2724	2728	2735	rBV4	45703	72598	1.19%	0.104%
67	19.198	2735	2739	2743	rBV	64934	96033	1.57%	0.138%
68	19.345	2760	2764	2769	rBV	127525	167200	2.74%	0.240%
69	19.551	2794	2799	2807	rBV	4208796	6101269	100.00%	8.761%
70	19.704	2821	2825	2828	rBV	125957	151035	2.48%	0.217%
71	19.745	2828	2832	2840	rVB7	50215	101259	1.66%	0.145%
72	20.080	2886	2889	2894	rVB4	90216	112914	1.85%	0.162%
73	20.245	2912	2917	2928	rBV3	396382	926846	15.19%	1.331%
74	20.451	2947	2952	2962	rBV2	603370	937268	15.36%	1.346%
75	20.528	2962	2965	2967	rVV2	92188	122119	2.00%	0.175%
76	20.563	2967	2971	2987	rVB3	533650	1136958	18.63%	1.633%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
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Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Title : SVOA CALIBRATION

77	20.763	2997	3005	3016	rBV	989975	1655541	27.13%	2.377%
78	20.851	3017	3020	3028	rVB7	126076	213877	3.51%	0.307%
79	21.033	3046	3051	3060	rBV3	240845	473057	7.75%	0.679%
80	21.304	3092	3097	3103	rBV	2924229	3713986	60.87%	5.333%
81	23.521	3467	3474	3481	rBV	3262883	5957956	97.65%	8.555%
82	23.674	3493	3500	3511	rVB	1954289	3882550	63.64%	5.575%

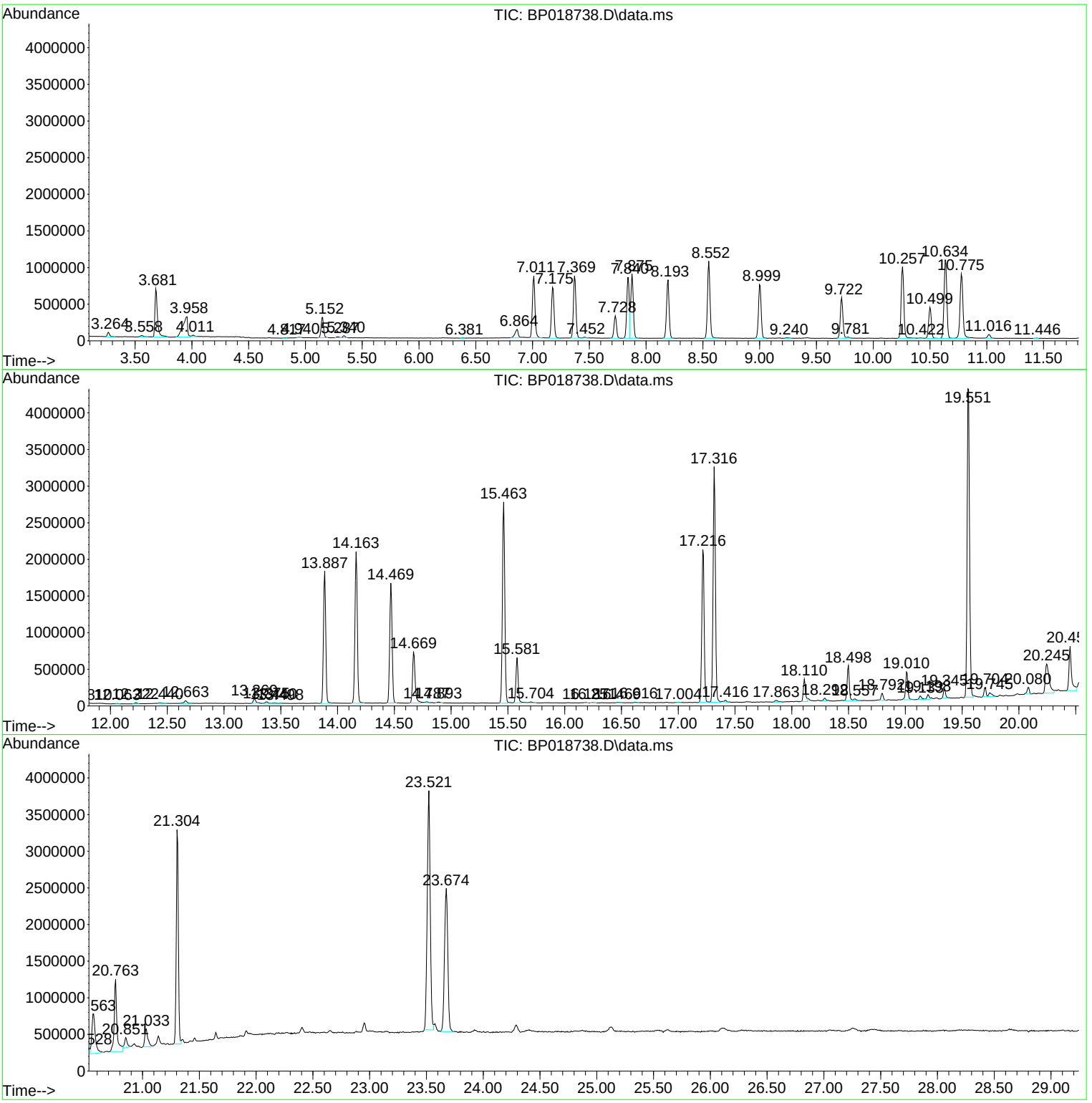
Sum of corrected areas: 69641362

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

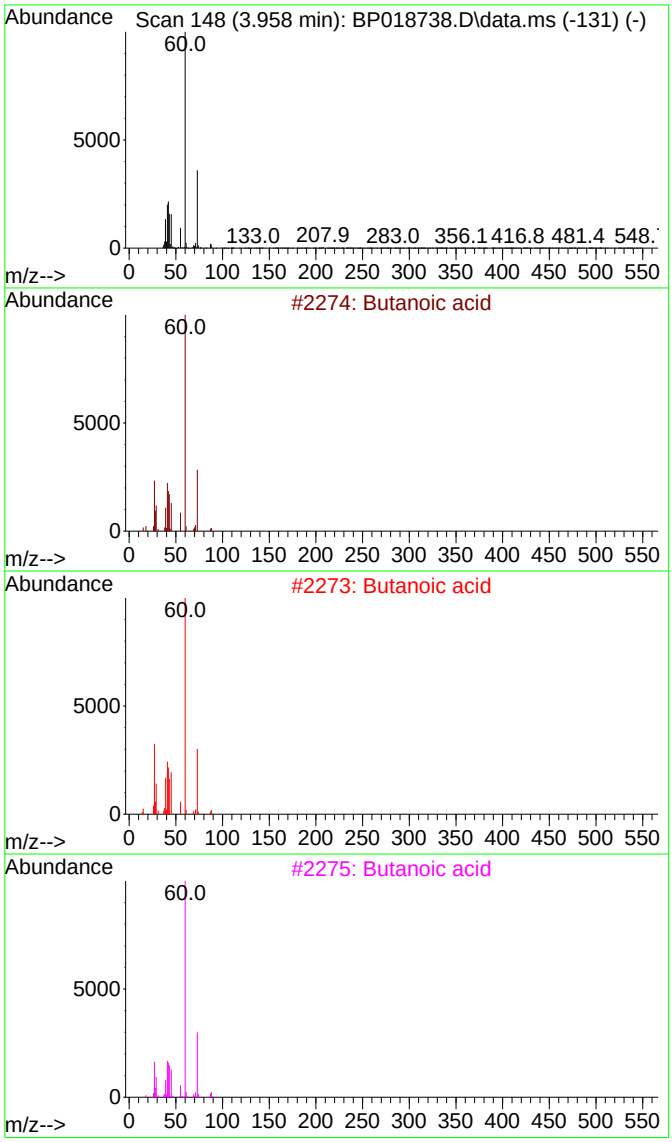
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

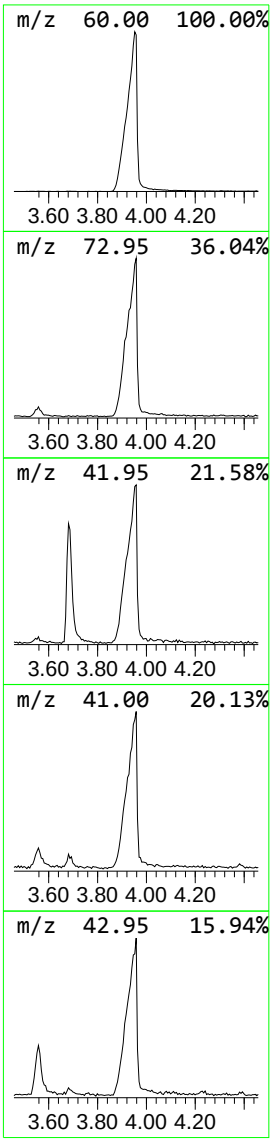
****

Peak Number 1 Butanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.958	12.35 ng/ul	824290	1,4-Dichlorobenzene-d4	7.840

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	91
2	Butanoic acid	88	C4H8O2	000107-92-6	91
3	Butanoic acid	88	C4H8O2	000107-92-6	80
4	Butanoic acid	88	C4H8O2	000107-92-6	72
5	Silver butanoate	194	C4H7AgO2	005076-24-4	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

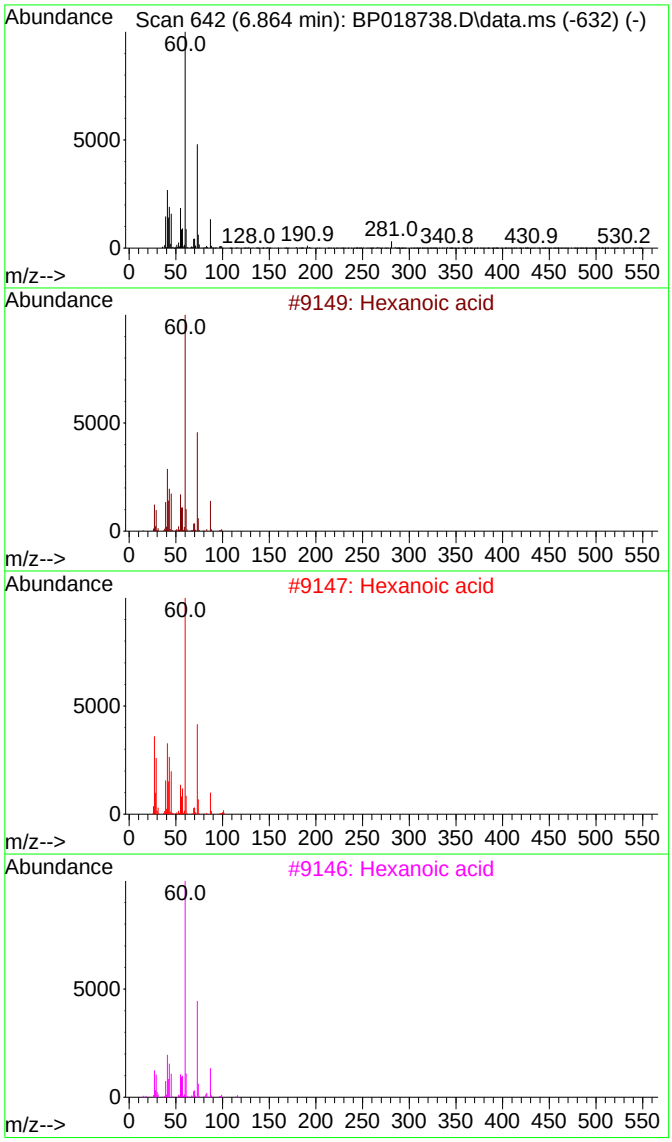
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

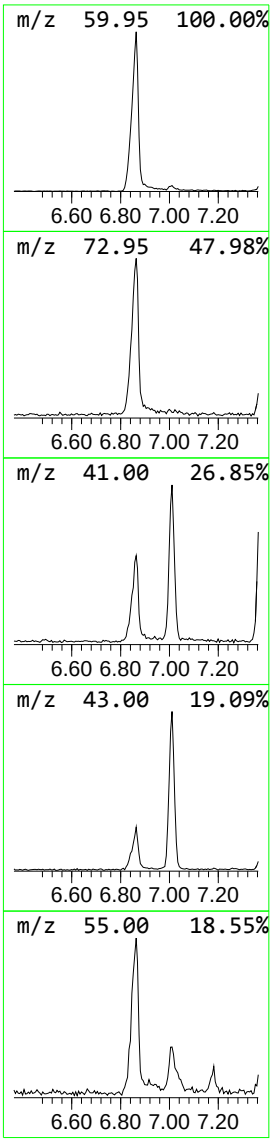
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.864	3.44 ng/ul	229819	1,4-Dichlorobenzene-d4	7.840

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid	116	C6H12O2	000142-62-1	90
2	Hexanoic acid	116	C6H12O2	000142-62-1	83
3	Hexanoic acid	116	C6H12O2	000142-62-1	83
4	Hexanoic acid	116	C6H12O2	000142-62-1	83
5	Pentanoic acid	102	C5H10O2	000109-52-4	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

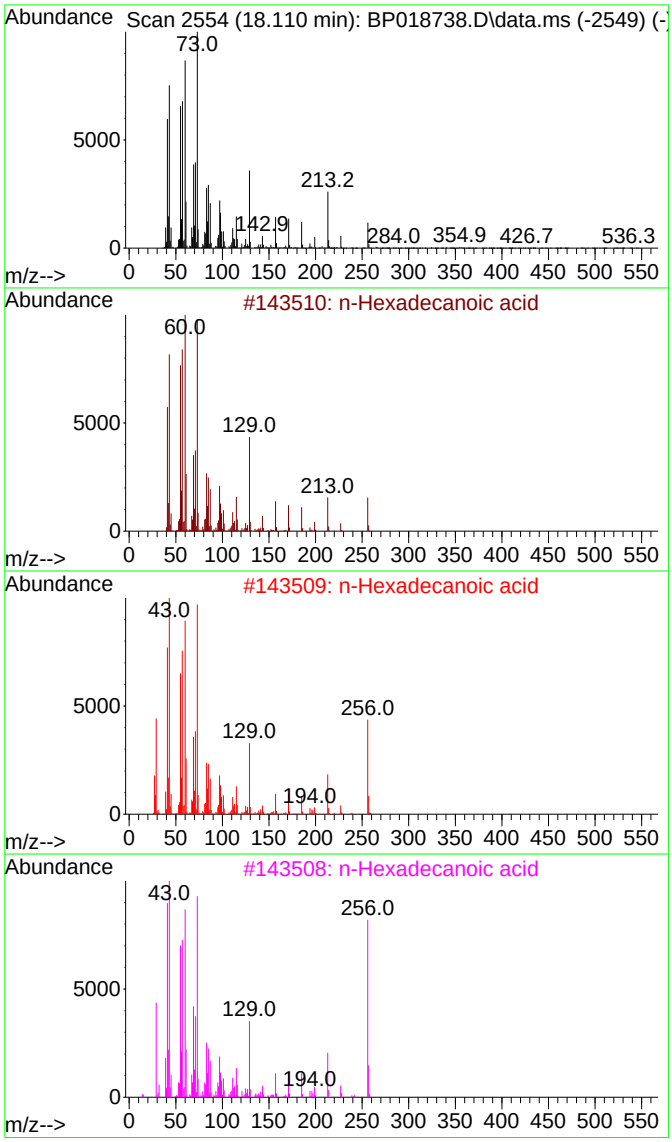
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

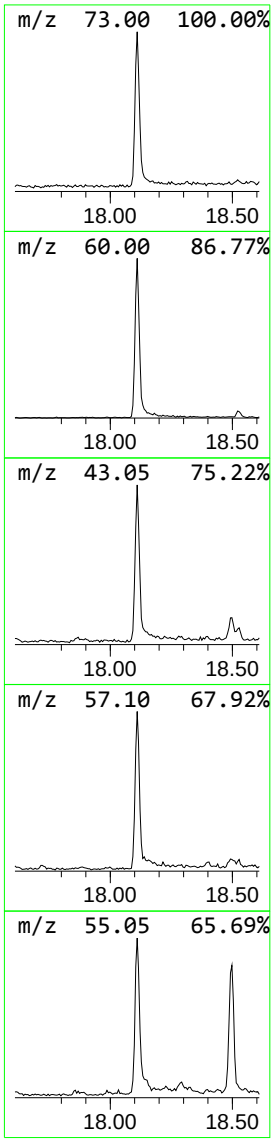
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.110	2.74 ng/ul	407968	Phenanthrene-d10	17.216

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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	Tridecanoic acid	214	C13H26O2	000638-53-9	91
5	Octadecanoic acid	284	C18H36O2	000057-11-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

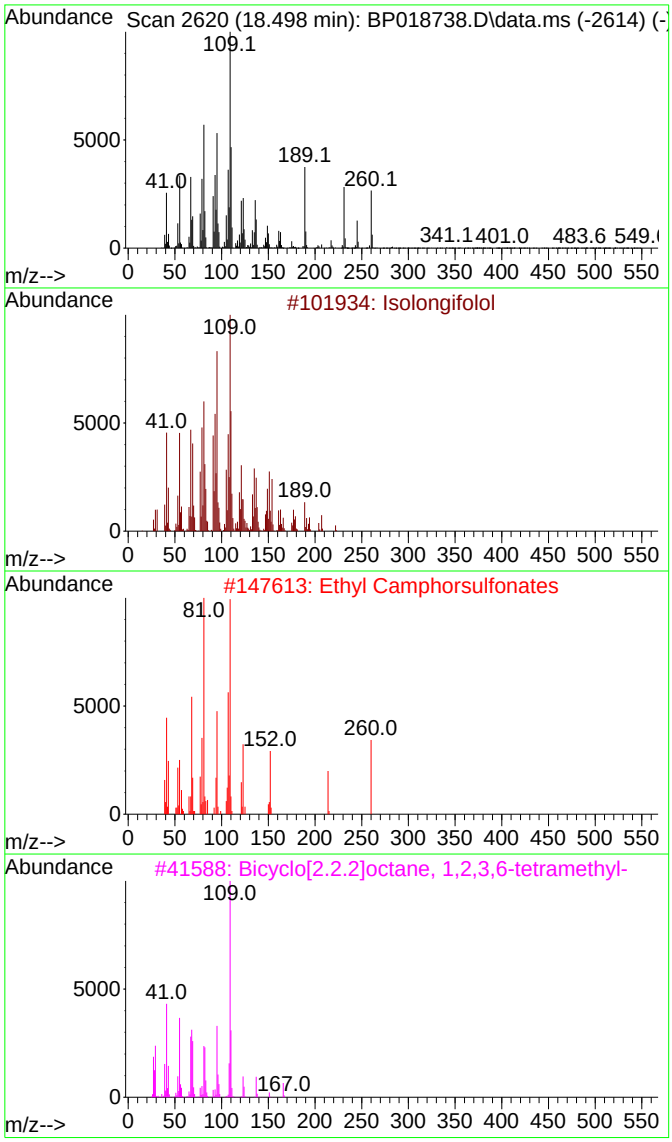
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

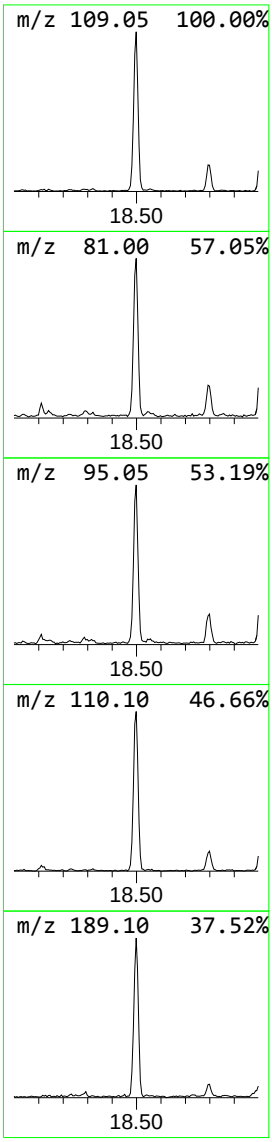
****

Peak Number 8 Isolongifolol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.498	4.33 ng/ul	644397	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isolongifolol	222	C15H26O	001139-17-9	64
2	Ethyl Camphorsulfonates	260	C12H20O4S	1000373-95-5	43
3	Bicyclo[2.2.2]octane, 1,2,3,6-te...	166	C12H22	062338-45-8	35
4	3,4,5-Trimethylpyrazole	110	C6H10N2	005519-42-6	27
5	2-Bornanecarboxylic acid	182	C11H18O2	1000163-82-4	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

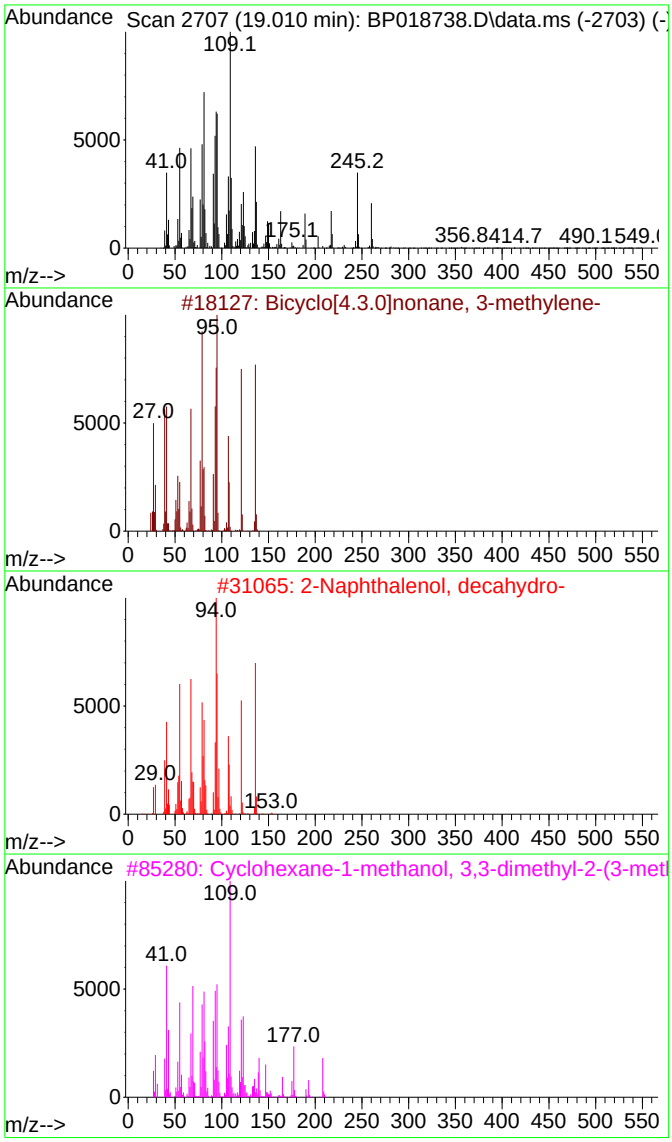
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

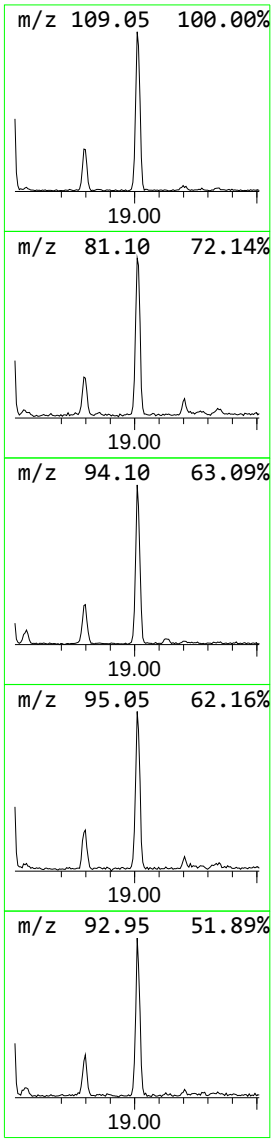
****

Peak Number 9 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.010	3.45 ng/ul	513177	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.3.0]nonane, 3-methylene-	136	C10H16	1000152-00-6	46
2	2-Naphthalenol, decahydro-	154	C10H18O	000825-51-4	38
3	Cyclohexane-1-methanol, 3,3-dime...	208	C14H24O	1000196-01-5	38
4	Isolongifolol	222	C15H26O	001139-17-9	35
5	Isoborneol,heptafluorobutyrate (...	350	C14H17F7O2	1000245-88-3	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

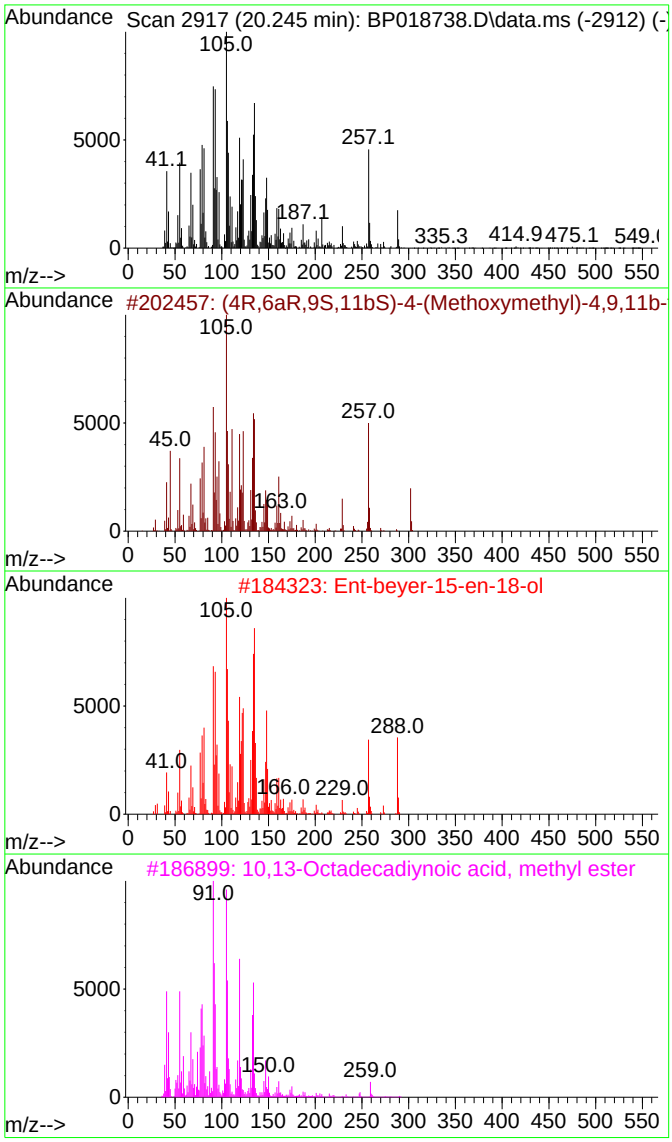
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

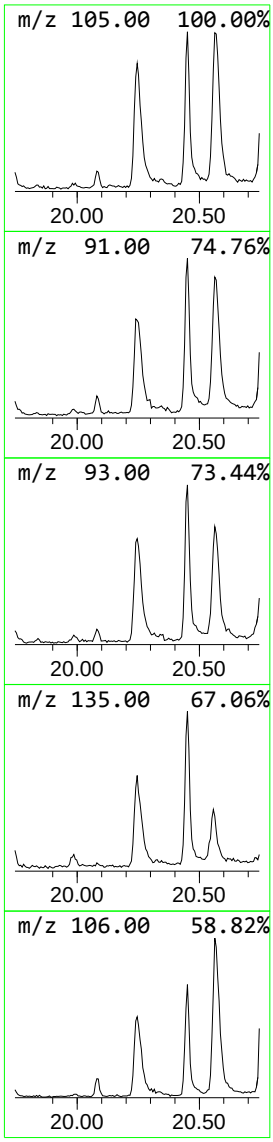
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (4R,6aR,9S,11bS)-4-(Methoxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.245	4.99 ng/ul	926846	Chrysene-d12	21.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(4R,6aR,9S,11bS)-4-(Methoxymethy...	302	C21H34O	1000495-78-0	52
2	Ent-beyer-15-en-18-ol	288	C20H32O	020107-90-8	38
3	10,13-Octadecadiynoic acid, meth...	290	C19H30O2	018202-24-9	38
4	Bicyclo[3.2.1]oct-2-ene, 3-methy...	134	C10H14	049826-53-1	30
5	6-Isopropenyl-3-methoxymethoxy-3...	196	C12H20O2	1000188-35-8	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

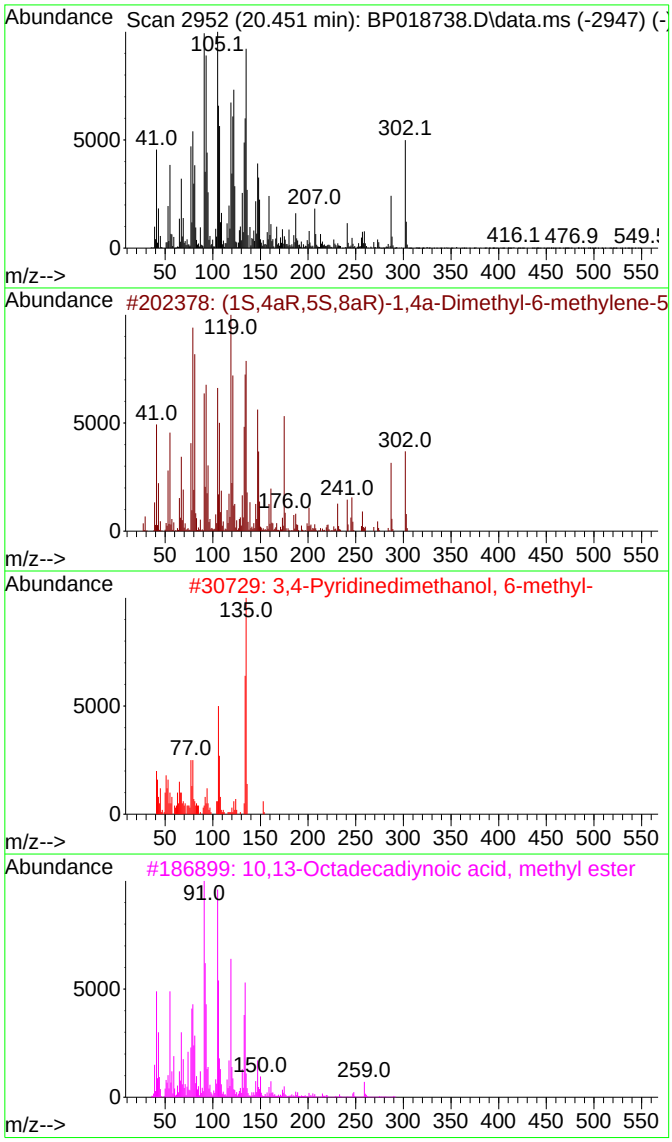
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

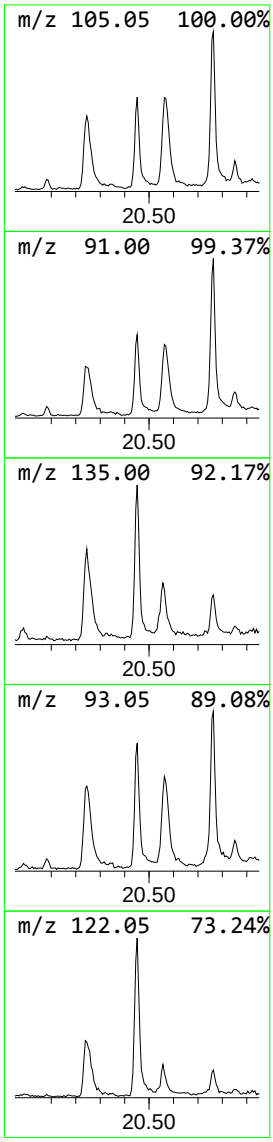
****

Peak Number 11 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.451	5.05 ng/ul	937268	Chrysene-d12	21.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S,4aR,5S,8aR)-1,4a-Dimethyl-6-...	302	C20H30O2	002761-77-5	38
2	3,4-Pyridinedimethanol, 6-methyl-	153	C8H11NO2	004664-11-3	30
3	10,13-Octadecadiynoic acid, meth...	290	C19H30O2	018202-24-9	30
4	Pyridine, 2,6-diethyl-	135	C9H13N	000935-28-4	27
5	6,7-Dehydrohyoscyamine	287	C17H21NO3	061616-97-5	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

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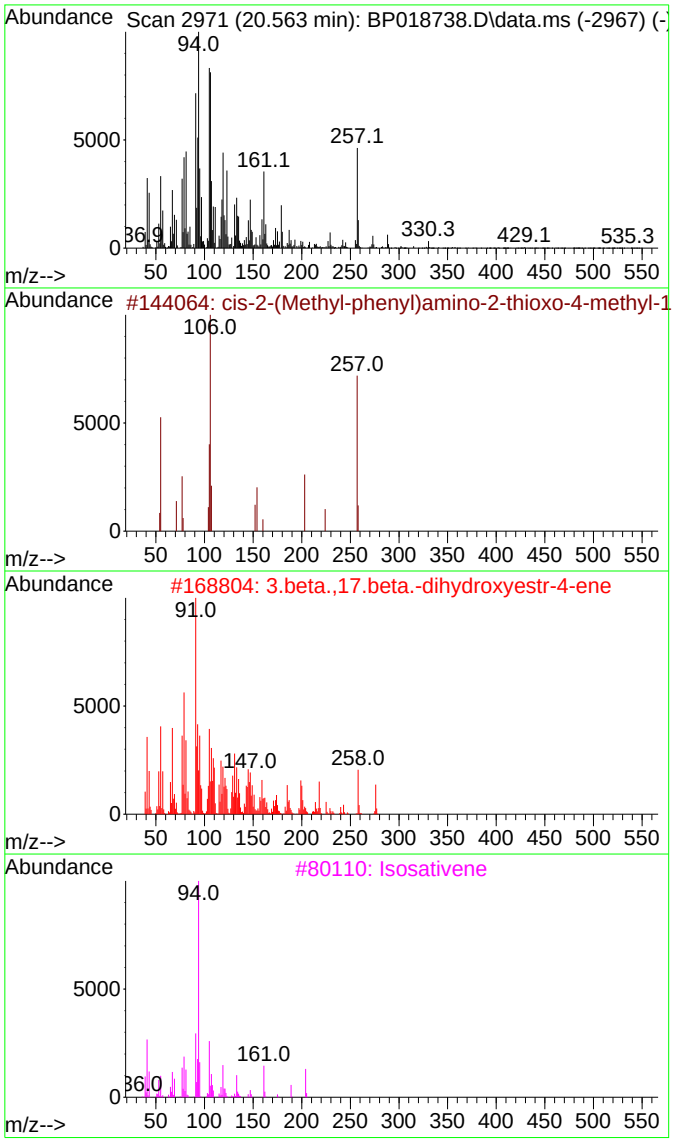
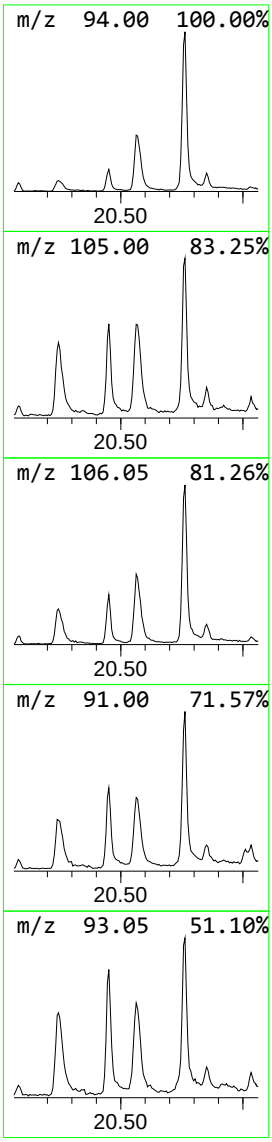
Peak Number 12 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.563	6.12 ng/ul	1136960	Chrysene-d12	21.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-2-(Methyl-phenyl)amino-2-thi...	257	C11H16NO2PS	1000145-42-9	25
2	3.beta.,17.beta.-dihydroxyestr-4...	276	C18H28O2	019793-20-5	11
3	Isosativene	204	C15H24	024959-83-9	11
4	2-Phenyl-hex-5-en-3-ol	176	C12H16O	077383-06-3	11
5	Amantadine	151	C10H17N	000768-94-5	11





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

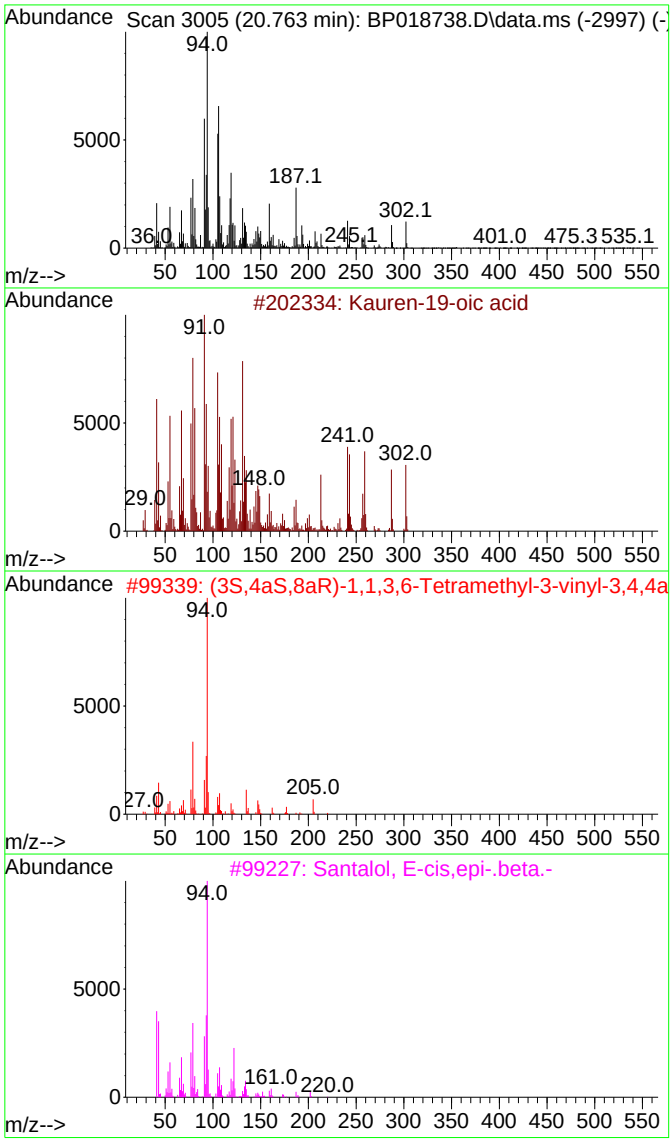
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

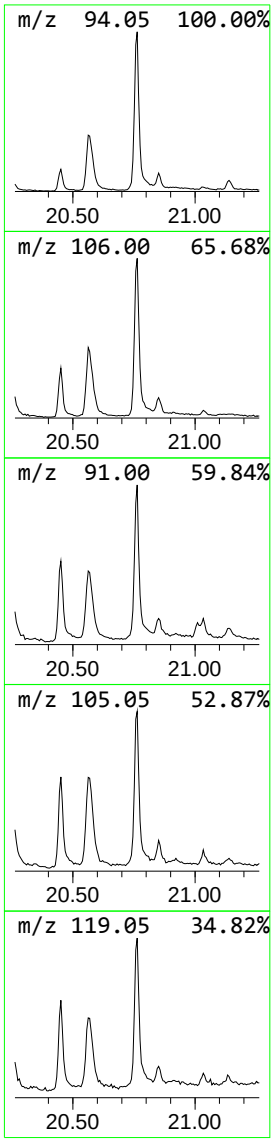
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Kauren-19-oic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.763	8.92 ng/ul	1655540	Chrysene-d12	21.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Kauren-19-oic acid	302	C20H30O2	006730-83-2	90
2	(3S,4aS,8aR)-1,1,3,6-Tetramethyl...	220	C15H24O	107602-53-9	27
3	Santalol, E-cis,epi-.beta.-	220	C15H24O	014490-17-6	27
4	(Z)-epi-.beta.-Santalol	220	C15H24O	079081-90-6	27
5	Menth-1-en-9-yl tiglate, p-	236	C15H24O2	1000383-62-5	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

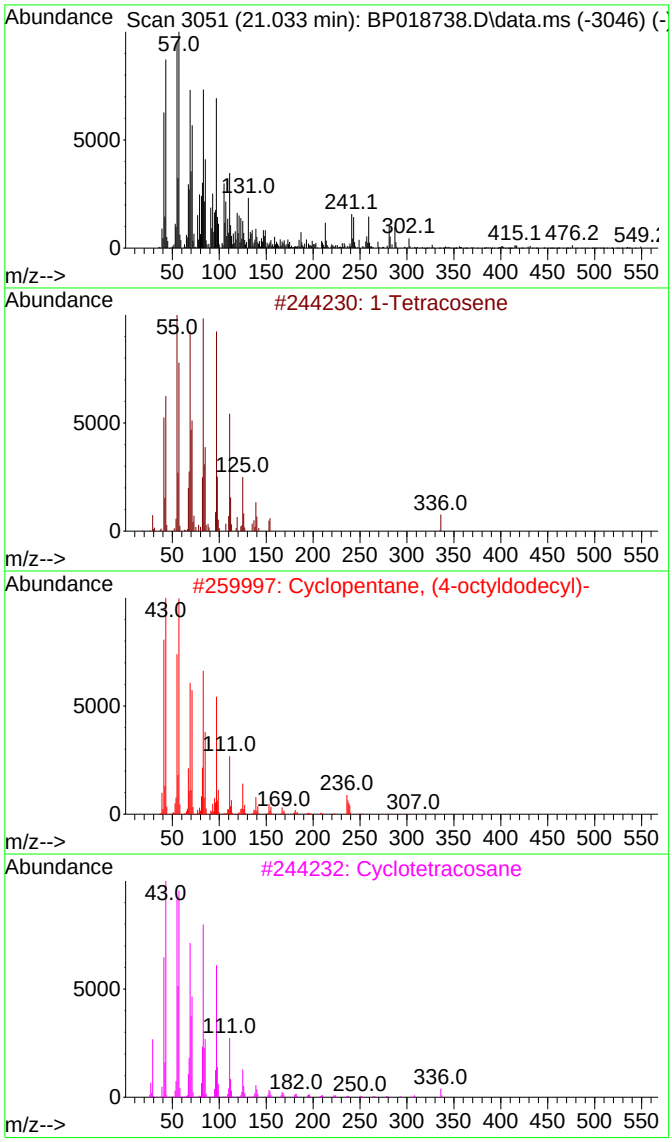
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

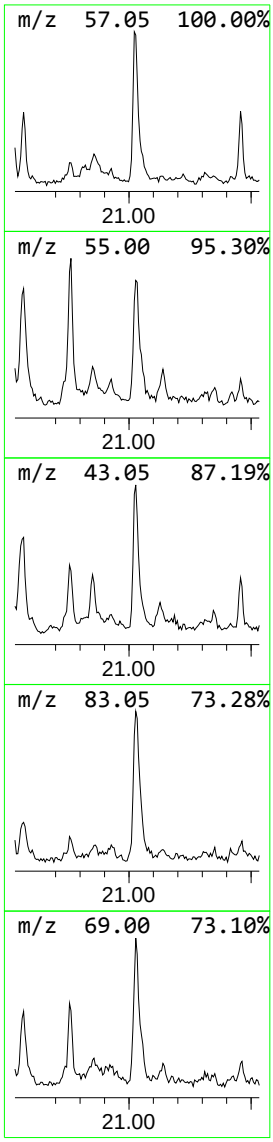
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.033	2.55 ng/ul	473057	Chrysene-d12	21.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	97
2	Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	92
3	Cyclotetracosane	336	C24H48	000297-03-0	90
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	90
5	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018738.D
Acq On : 11 Dec 2023 21:41
Operator : MA/JU
Sample : 04929-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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
Butanoic acid	3.958	12.4	ng/ul	824290	1	7.840	1335070	20.0
Hexanoic acid	6.864	3.4	ng/ul	229819	1	7.840	1335070	20.0
n-Hexadecanoic ...	18.110	2.7	ng/ul	407968	4	17.216	2977860	20.0
Isolongifolol	18.498	4.3	ng/ul	644397	4	17.216	2977860	20.0
unknown-01	19.010	3.5	ng/ul	513177	4	17.216	2977860	20.0
(4R,6aR,9S,11bS...	20.245	5.0	ng/ul	926846	5	21.304	3713990	20.0
unknown-02	20.451	5.0	ng/ul	937268	5	21.304	3713990	20.0
unknown-03	20.563	6.1	ng/ul	1136960	5	21.304	3713990	20.0
Kauren-19-oic acid	20.763	8.9	ng/ul	1655540	5	21.304	3713990	20.0
(DEL) Alkane: S...	21.033	2.5	ng/ul	473057	5	21.304	3713990	20.0