LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015090.D
Acq On : 25 Jun 2021 22:31
Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGD44

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

Signal : TIC: BN015090.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	31	34	42	rVB	28577	35238	1.48%	0.104%
2	3.540	73	77	93	rBV	303744	483007	20.29%	1.431%
3	3.681	98	101	112	rBV	218783	324821	13.64%	0.963%
4	3.775	113	117	124	rVB3	24186	42821	1.80%	0.127%
5	3.875	128	134	139	rBV	62525	96878	4.07%	0.287%
6	4.005	151	156	167	rVB	58843	95022	3.99%	0.282%
7	4.358	211	216	224	rVB3	25195	38180	1.60%	0.113%
8	4.487	234	238	242	rBV	18021	23447	0.98%	0.069%
9	4.534	242	246	251	rBV	52248	71719	3.01%	0.213%
10	4.887	303	306	311	rVB2	8207	13990	0.59%	0.041%
11	5.228	354	364	369	rBV	95249	215831	9.06%	0.640%
12	5.281	369	373	380	rVB	44342	67095	2.82%	0.199%
13	5.411	390	395	404	rVB	124247	250928	10.54%	0.744%
14	5.775	451	457	462	rBV	63941	92870	3.90%	0.275%
15	5.840	466	468	476	rVB3	5504	9191	0.39%	0.027%
16	6.722	612	618	627	rBV4	18244	45129	1.90%	0.134%
17	6.893	640	647	656	rBV	387854	618418	25.97%	1.833%
18	6.981	657	662	667	rVB7	4292	8892	0.37%	0.026%
19	7.063	667	676	685	rBV	309794	492493	20.68%	1.460%
20	7.263	703	710	720	rVB	391643	633707	26.62%	1.878%
21	7.363	720	727	730	rBV6	7124	12150	0.51%	0.036%
22	7.728	783	789	796	rVV	607089	951146	39.95%	2.819%
23	7.787	796	799	807	rVB4	5860	10376	0.44%	0.031%
24	7.869	807	813	822	rVB6	5059	13031	0.55%	0.039%
25	7.963	825	829	837	rVB2	11392	19958	0.84%	0.059%
26	8.275	876	882	895	rVB3	13259	27658	1.16%	0.082%
27	8.416	899	906	914	rBV	453855	717744	30.15%	2.127%
28	8.534	921	926	933	rVB	20767	32650	1.37%	0.097%
29	8.816	968	974	977	rBV2	16334	25384	1.07%	0.075%
30	8.869	977	983	994	rVB	355393	572665	24.05%	1.697%
31	9.587	1098	1105	1112	rVV	207166	339527	14.26%	1.006%
32	10.116	1187	1195	1204	rBV	462208	745976	31.33%	2.211%
33	10.281	1216	1223	1233	rBV	43479	76516	3.21%	0.227%
34	10.499	1252	1260	1268	rBV	822144	1348391	56.63%	3.996%
35	10.628	1275	1282	1291	rBV	392857	667310	28.03%	1.978%
36	11.375	1405	1409	1419	rBV	18947	33869	1.42%	0.100%

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Instrument : BNA_N
ClientSampleId : BGD44

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

37	12.087	1525	1530	1534	rVB	7304	11231	0.47%	0.033%
38	12.263	1556	1560	1565	rVV2	5589	9152	0.38%	0.027%
39	12.716	1631	1637	1641	rBV6	5013	9452	0.40%	0.028%
40	13.193	1713	1718	1722	rBV	16378	23145	0.97%	0.069%
41	13.763	1809	1815	1823	rBV	767998	1044719	43.88%	3.096%
42	14.040	1855	1862	1871	rVV	871495	1322237	55.53%	3.919%
43	14.269	1895	1901	1906	rBV	25321	32830	1.38%	0.097%
44	14.351	1908	1915	1921	rVV	1183600	1733140	72.79%	5.136%
45	14.392	1921	1922	1929	rVV2	17301	16383	0.69%	0.049%
46	14.469	1929	1935	1938	rVV4	6553	9102	0.38%	0.027%
47	14.539	1941	1947	1958	rVV	151860	233543	9.81%	0.692%
48	14.698	1968	1974	1980	rBV7	14528	25840	1.09%	0.077%
49	14.775	1982	1987	1992	rVB4	14251	17746	0.75%	0.053%
50	15.181	2051	2056	2059	rBV3	6611	10356	0.43%	0.031%
51	15.222	2059	2063	2069	rVB	17388	23044	0.97%	0.068%
52	15.345	2077	2084	2091	rBV	1175517	1634888	68.67%	4.845%
53	15.457	2098	2103	2109	rVV2	35031	49712	2.09%	0.147%
54	15.516	2109	2113	2116	rVB6	7500	9654	0.41%	0.029%
55	15.586	2121	2125	2129	rVV2	11555	17256	0.72%	0.051%
56	15.775	2152	2157	2164	rVV	239564	300138	12.61%	0.889%
57	16.045	2198	2203	2206	rBV5	5840	10932	0.46%	0.032%
58	16.086	2206	2210	2211	rBV2	9545	12085	0.51%	0.036%
59	16.369	2252	2258	2259	rBV6	8516	14886	0.63%	0.044%
60	16.404	2260	2264	2268	rVV	23410	30176	1.27%	0.089%
61	16.492	2275	2279	2285	rBV8	11117	21708	0.91%	0.064%
62	16.628	2299	2302	2305	rVB4	7880	9828	0.41%	0.029%
63	16.739	2318	2321	2325	rVB5	10855	12106	0.51%	0.036%
64	16.881	2341	2345	2348	rBV2	26829	34562	1.45%	0.102%
65	16.922	2348	2352	2357	rVV6	13082	23225	0.98%	0.069%
66	16.963	2357	2359	2362	rVB4	7136	8941	0.38%	0.026%
67	17.033	2367	2371	2374	rVB5	9466	12576	0.53%	0.037%
68	17.098	2374	2382	2387	rBV	1456336	2047385	85.99%	6.068%
69	17.139	2387	2389	2393	rVV	59036	76072	3.20%	0.225%
70	17.198	2393	2399	2407	rVB	1215683	1785335	74.98%	5.291%
71	17.345	2422	2424	2429	rBV5	11380	18460	0.78%	0.055%
72	17.410	2434	2435	2439	rBV4	7448	10343	0.43%	0.031%
73	17.469	2442	2445	2449	rBV6	16293	28953	1.22%	0.086%
74	17.545	2456	2458	2461	rVB4	19073	19438	0.82%	0.058%
75	17.616	2467	2470	2477	rVB	38578	46646	1.96%	0.138%
76	17.786	2495	2499	2505	rVB4	25226	36141	1.52%	0.107%

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Instrument : BNA_N
ClientSampleId : BGD44

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

77	17.975	2527	2531	2533	rBV2	41347	50433	2.12%	0.149%
78	18.004	2533	2536	2543	rVV6	25427	51692	2.17%	0.153%
79	18.169	2560	2564	2567	rBV3	16740	27033	1.14%	0.080%
80	18.251	2574	2578	2583	rVB2	30342	43478	1.83%	0.129%
81	18.310	2584	2588	2592	rBV2	47715	57861	2.43%	0.171%
82	18.428	2604	2608	2612	rVB	46893	65171	2.74%	0.193%
83	18.475	2613	2616	2620	rVB6	22813	31239	1.31%	0.093%
84	18.945	2692	2696	2702	rBV4	69487	133807	5.62%	0.397%
85	19.163	2729	2733	2740	rVB	165403	239119	10.04%	0.709%
86	19.498	2784	2790	2794	rBV	1501648	2229038	93.62%	6.606%
87	20.063	2883	2886	2890	rBV	37925	43698	1.84%	0.130%
88	20.516	2958	2963	2968	rBV	1203913	1381343	58.02%	4.094%
89	21.027	3047	3050	3053	rBV3	60052	61548	2.59%	0.182%
90	21.221	3079	3083	3089	rVB	339247	377548	15.86%	1.119%
91	21.292	3089	3095	3100	rBV	1782783	2353709	98.86%	6.975%
92	22.157	3238	3242	3246	rBV	234629	308050	12.94%	0.913%
93	22.363	3275	3277	3286	rVB2	126848	195691	8.22%	0.580%
94	22.874	3360	3364	3374	rVB3	197223	399190	16.77%	1.183%
95	23.398	3446	3453	3458	rBV	1093449	2014402	84.60%	5.970%
96	23.451	3458	3462	3469	rVB2	208583	359874	15.11%	1.067%
97	23.539	3470	3477	3490	rVB2	1275634	2380956	100.00%	7.056%
98	24.098	3567	3572	3580	rVB	176201	351608	14.77%	1.042%
99	24.857	3695	3701	3712	rVB	158618	372727	15.65%	1.105%
100	25.739	3845	3851	3859	rVB2	74878	166584	7.00%	0.494%

Sum of corrected areas: 33743193

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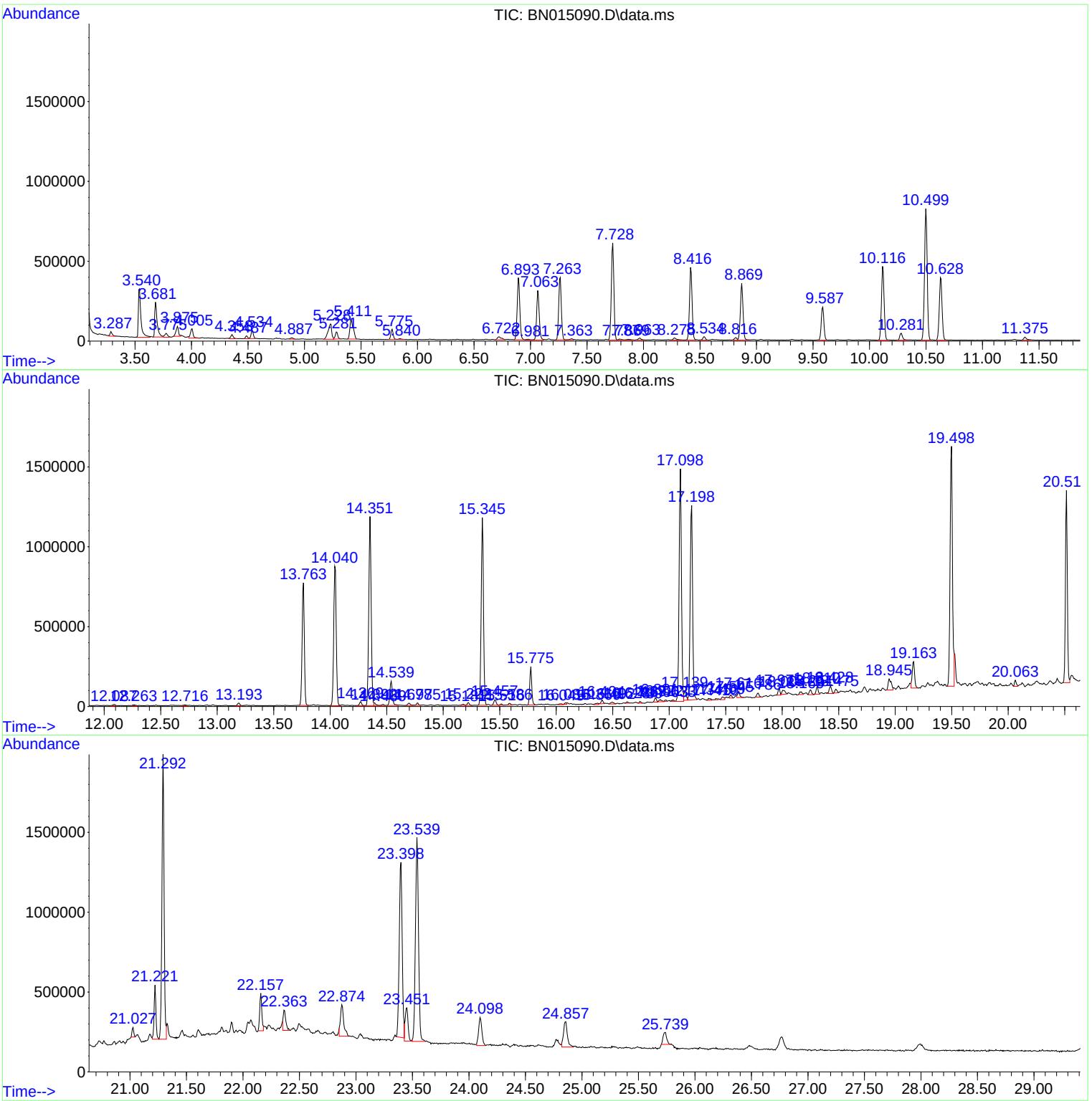
Instrument :
BNA_N
ClientSampleId :
BGD44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Quant Title : SVOA CALIBRATION

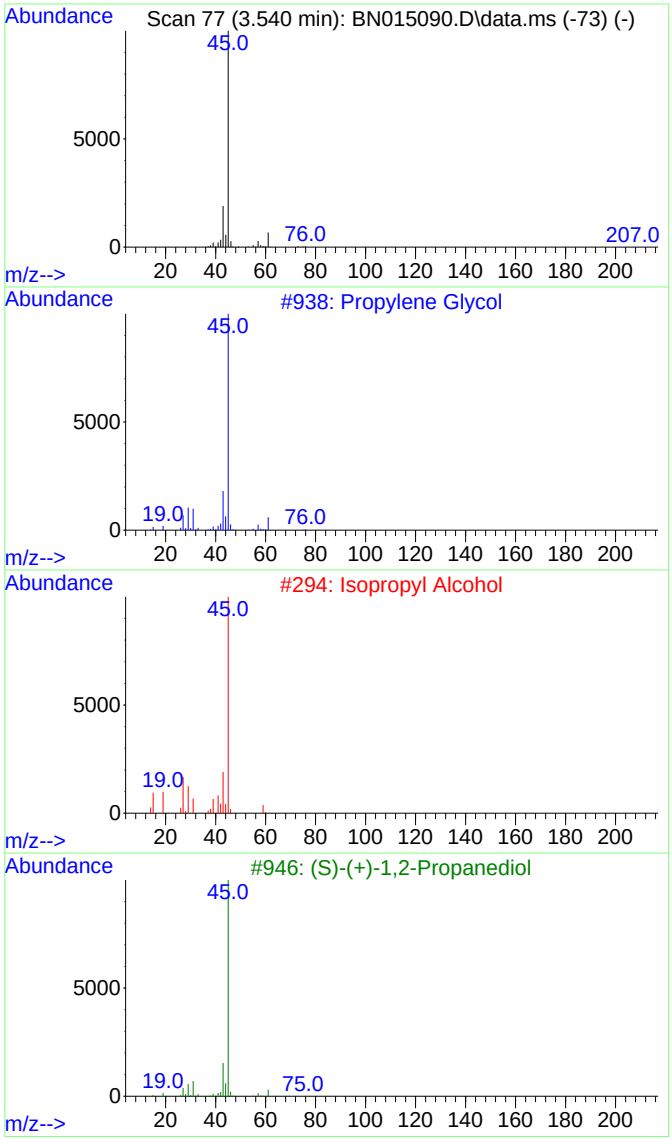
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

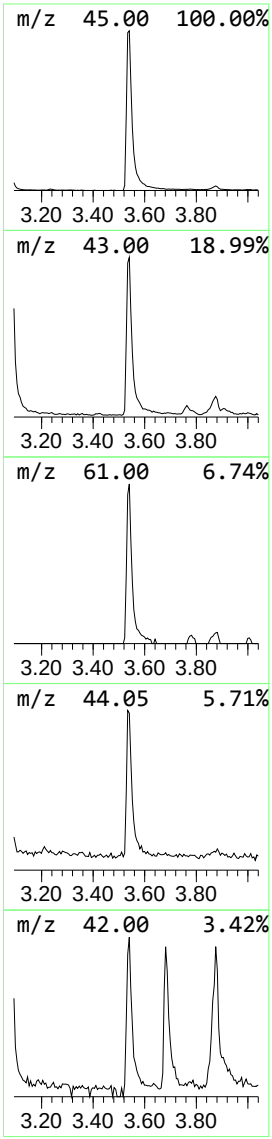
****

Peak Number 1 Propylene Glycol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	10.16 ng/ul	483007	1,4-Dichlorobenzene-d4	7.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	90
2	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	37
5	Isopropyl Alcohol	60	C3H8O	000067-63-0	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

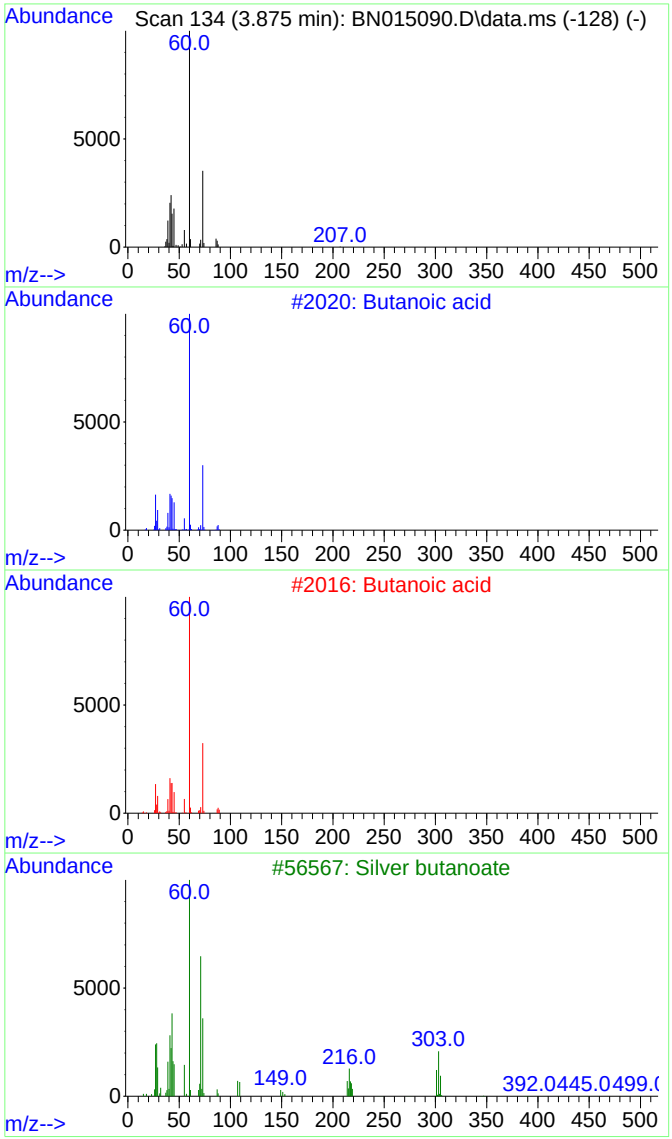
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

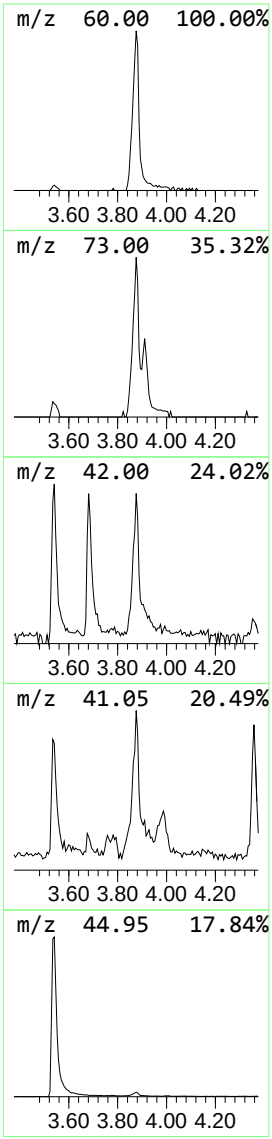
****

Peak Number 2 Butanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.875	2.04 ng/ul	96878	1,4-Dichlorobenzene-d4	7.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	72
2	Butanoic acid	88	C4H8O2	000107-92-6	72
3	Silver butanoate	194	C4H7AgO2	005076-24-4	64
4	.beta.-l-Arabinopyranoside, methyl	164	C6H12O5	001825-00-9	40
5	Pentanoic acid	102	C5H10O2	000109-52-4	39





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

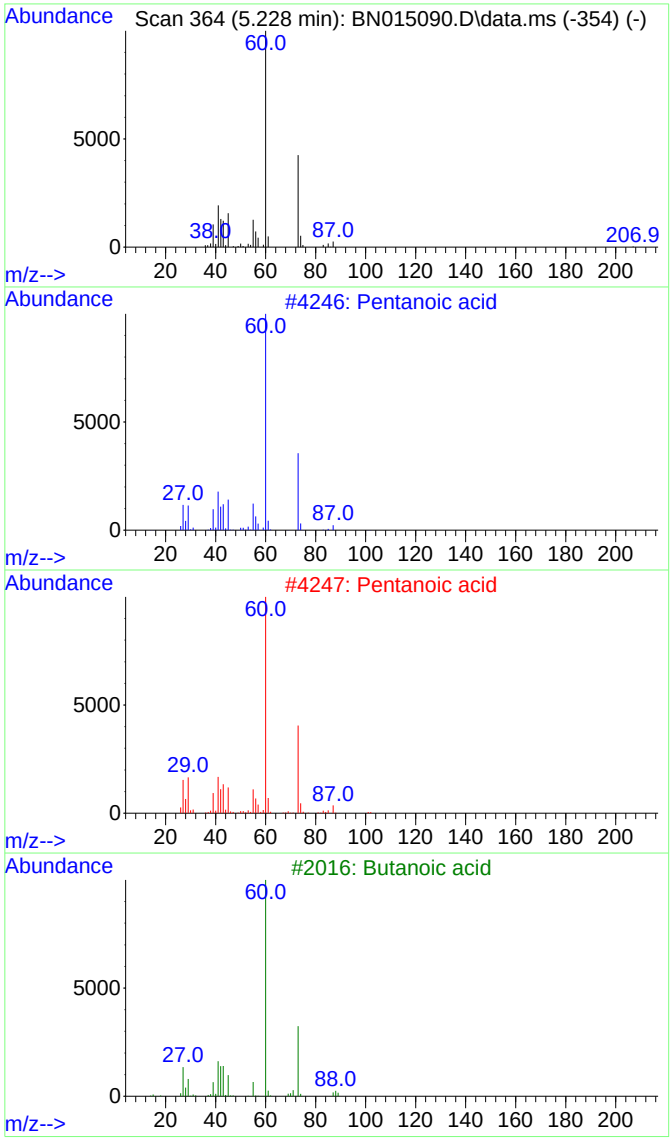
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

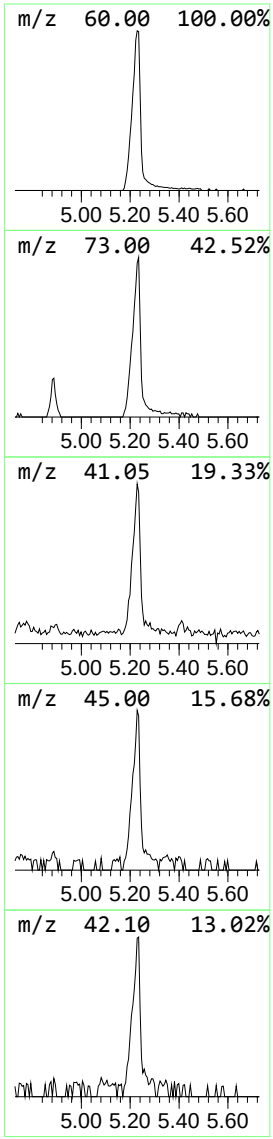
****

Peak Number 3 Pentanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.228	4.54 ng/ul	215831	1,4-Dichlorobenzene-d4	7.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid	102	C5H10O2	000109-52-4	90
2	Pentanoic acid	102	C5H10O2	000109-52-4	83
3	Butanoic acid	88	C4H8O2	000107-92-6	78
4	Pentanoic acid	102	C5H10O2	000109-52-4	72
5	Butanoic acid	88	C4H8O2	000107-92-6	64





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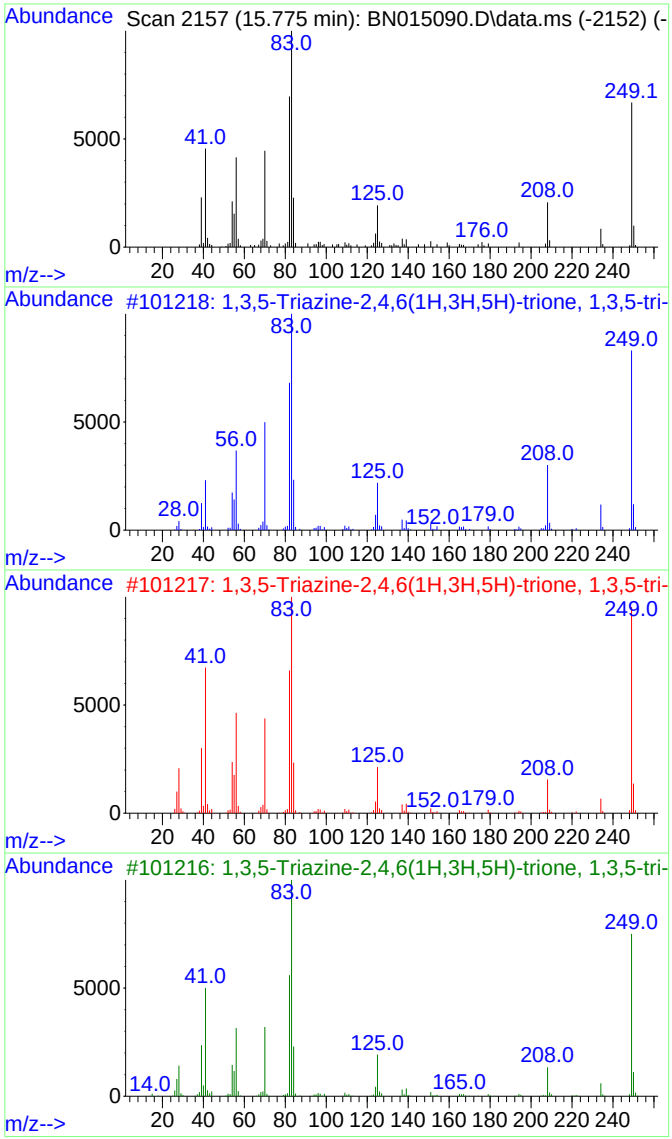
Quant Title : SVOA CALIBRATION

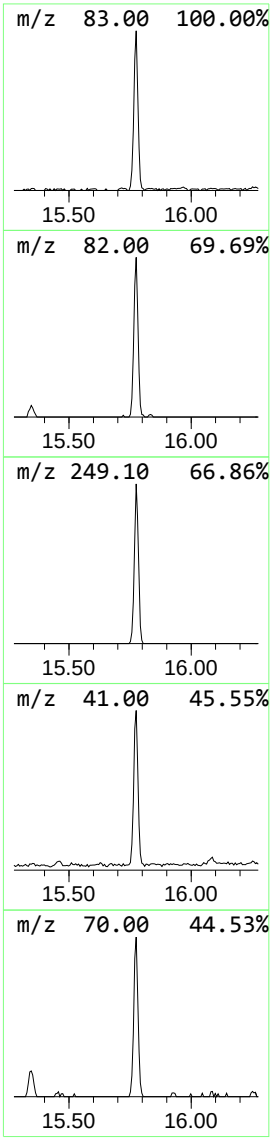
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.93 ng/ul	300138	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	99
2	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	96
3	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	93
4	4H-1-Benzopyran-2-carboxylic aci...	249	C12H11NO5	030192-51-9	23
5	4H-1-Benzopyran-2-carboxylic aci...	249	C12H11NO5	032142-43-1	23





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Instrument :
BNA_N
ClientSampleId :
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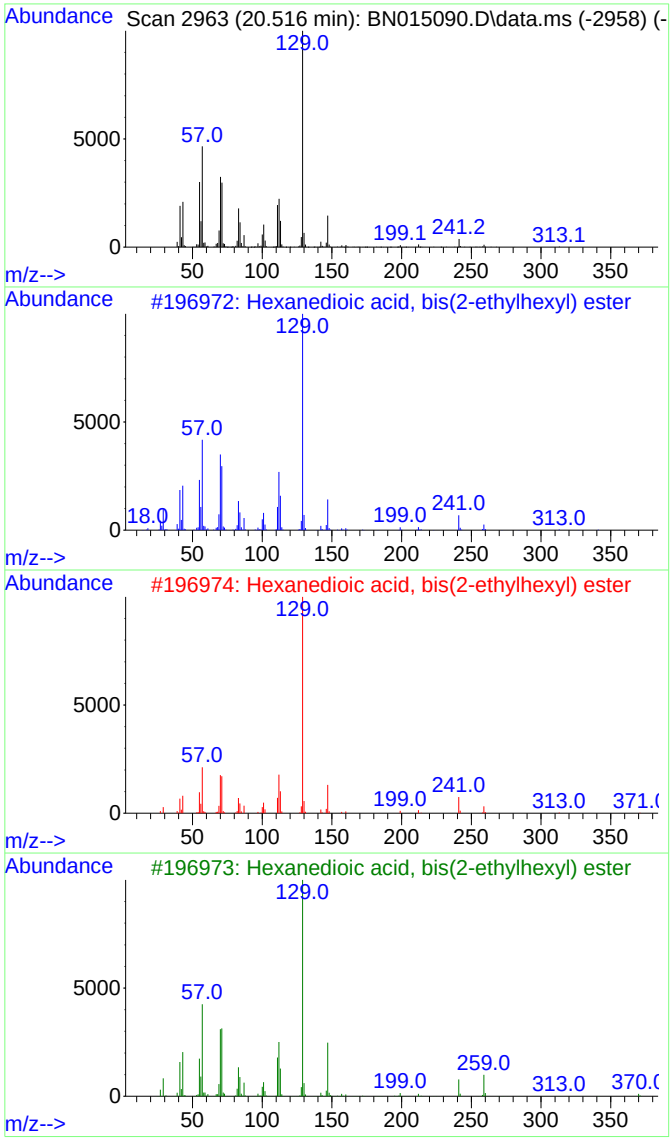
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

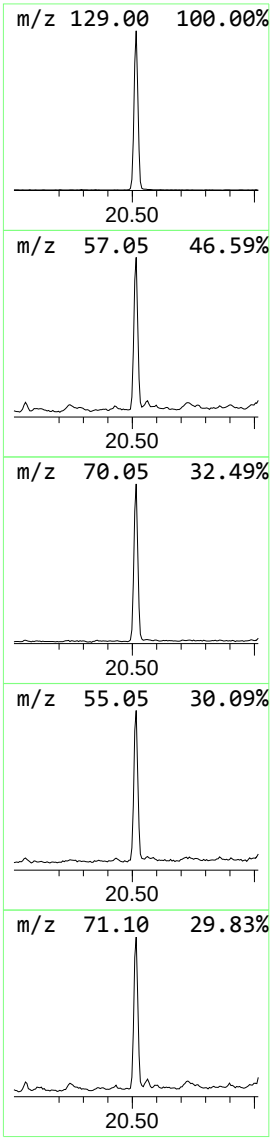
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	11.74 ng/ul	1381340	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76
5	Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	68





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Data File : BN015090.D
Acq On : 25 Jun 2021 22:31
Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGD44

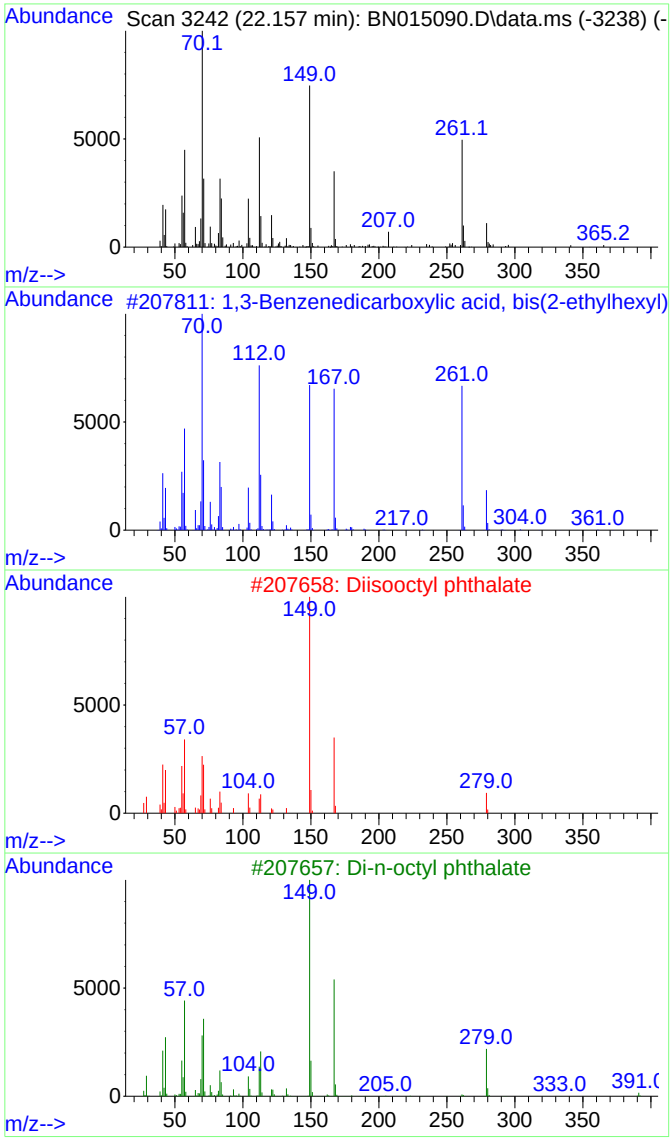
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

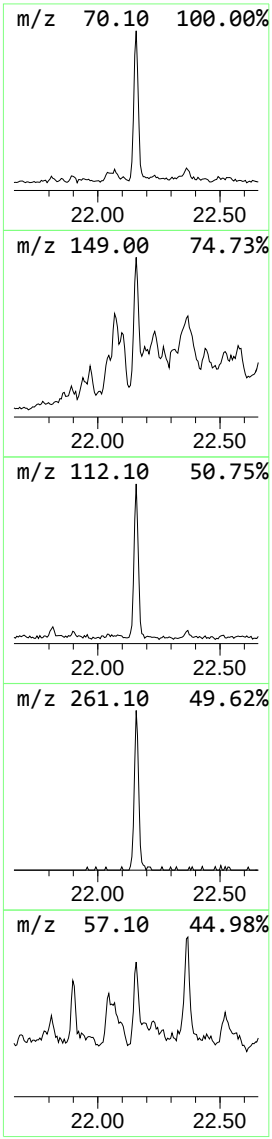
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,3-Benzenedicarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.157	2.62 ng/ul	308050	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	53
2	Diisooctyl phthalate	390	C24H38O4	000131-20-4	35
3	Di-n-octyl phthalate	390	C24H38O4	000117-84-0	27
4	Diisooctyl phthalate	390	C24H38O4	000131-20-4	27
5	9-(2',2'-Dimethylpropanoilhydraz...	576	C30H42Cl2N4O3	1000111-04-6	27





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Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD44

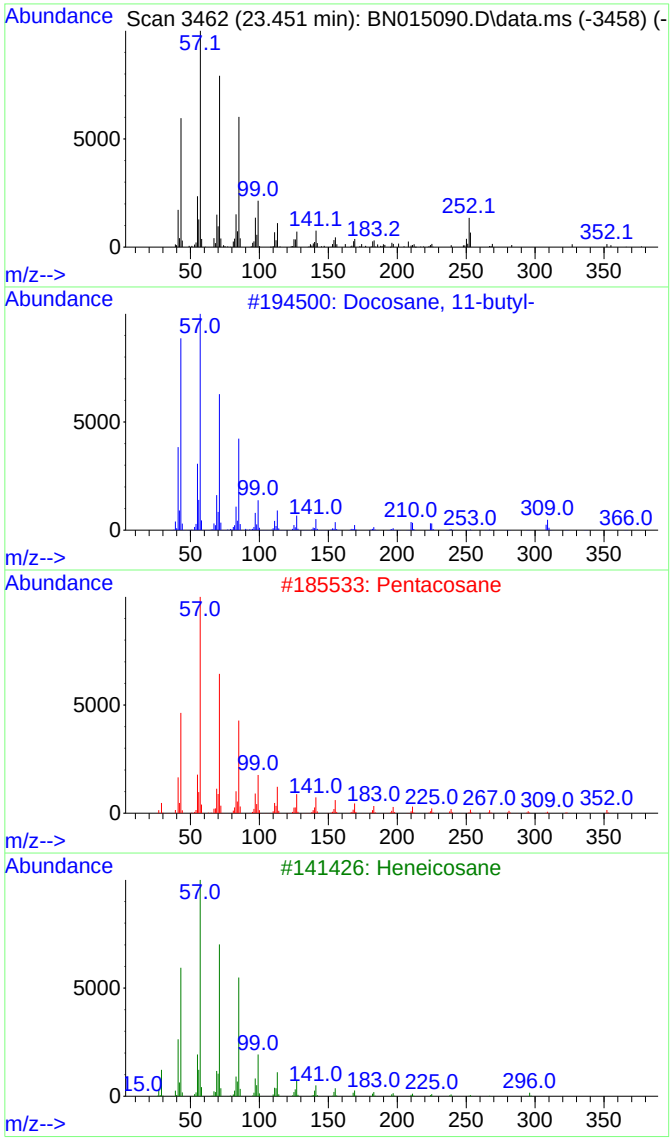
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

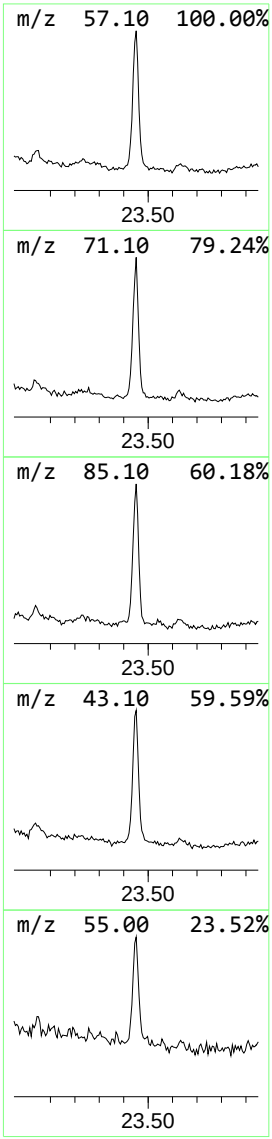
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.451	3.02 ng/ul	359874	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	87
5	Heneicosane	296	C21H44	000629-94-7	87





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Data File : BN015090.D
Acq On : 25 Jun 2021 22:31
Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD44

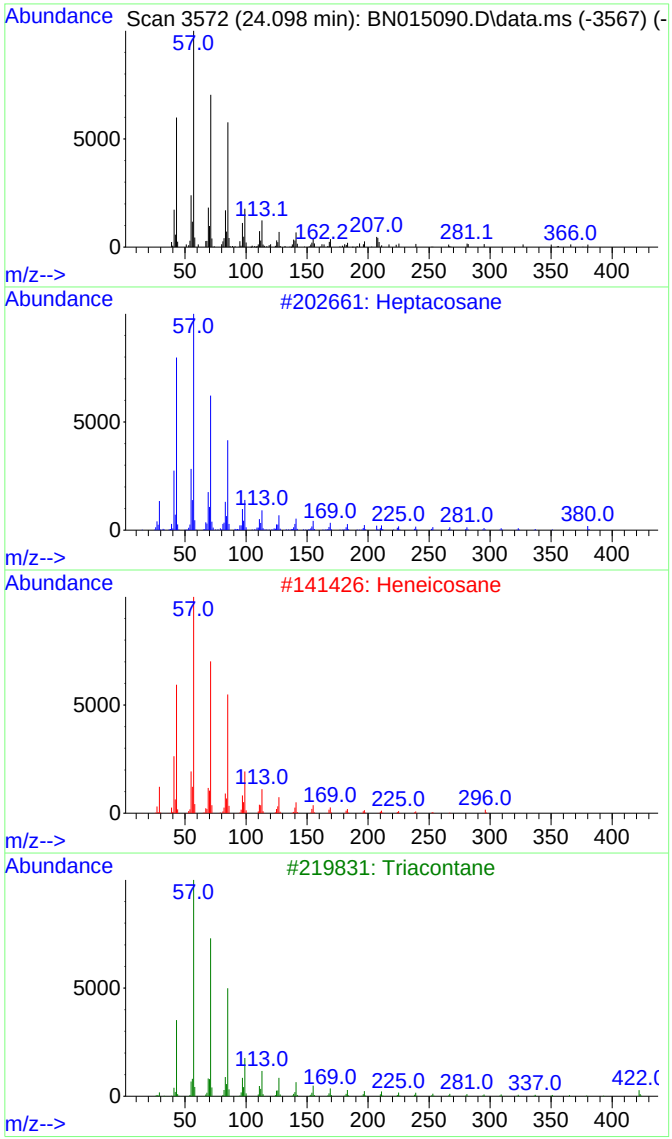
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

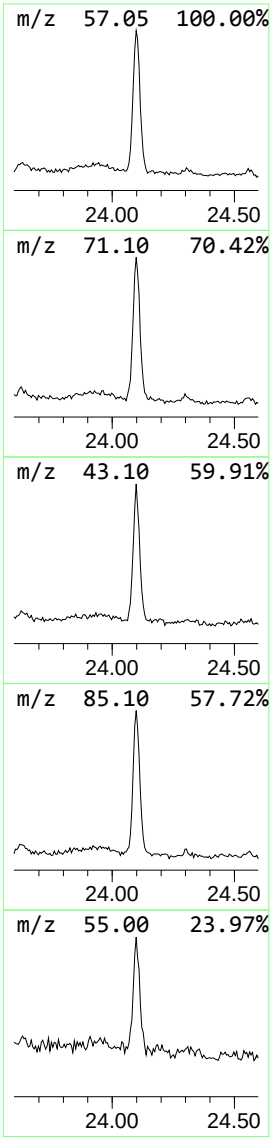
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.098	2.95 ng/ul	351608	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	95
2	Heneicosane	296	C21H44	000629-94-7	90
3	Triacontane	422	C30H62	000638-68-6	90
4	2-methyloctacosane	408	C29H60	1000376-72-8	90
5	Eicosane	282	C20H42	000112-95-8	90





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Data File : BN015090.D
Acq On : 25 Jun 2021 22:31
Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGD44

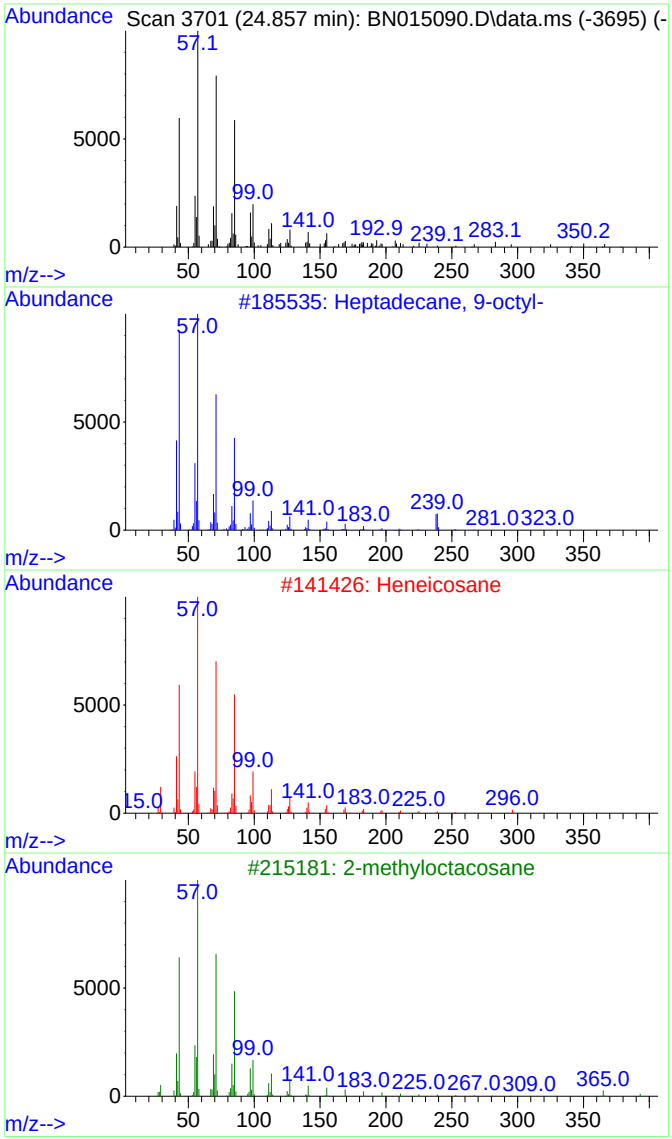
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

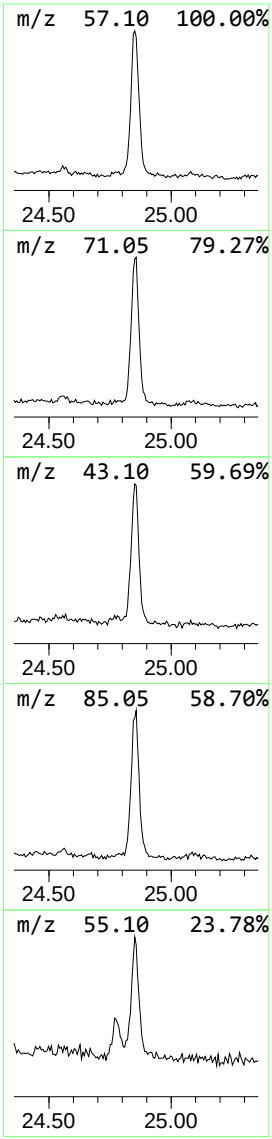
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.857	3.13 ng/ul	372727	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Octacosane	394	C28H58	000630-02-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015090.D
Acq On : 25 Jun 2021 22:31
Operator : CG/JU
Sample : M2680-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGD44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propylene Glycol	3.540	10.2	ng/ul	483007	1	7.728	951146	20.0
Butanoic acid	3.875	2.0	ng/ul	96878	1	7.728	951146	20.0
Pentanoic acid	5.228	4.5	ng/ul	215831	1	7.728	951146	20.0
1,3,5-Triazine-...	15.775	2.9	ng/ul	300138	4	17.098	2047390	20.0
Hexanedioic aci...	20.516	11.7	ng/ul	1381340	5	21.292	2353710	20.0
1,3-Benzenedica...	22.157	2.6	ng/ul	308050	5	21.292	2353710	20.0
(DEL) Alkane: S...	23.451	3.0	ng/ul	359874	6	23.539	2380960	20.0
(DEL) Alkane: S...	24.098	3.0	ng/ul	351608	6	23.539	2380960	20.0
(DEL) Alkane: S...	24.857	3.1	ng/ul	372727	6	23.539	2380960	20.0