LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071621\
Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

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_G\METHODS\SFAM-EPA-BG070921.M
Title : SVOA CALIBRATION

Signal : TIC: BG049389.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.102	5	11	20	rBV2	52915	132798	9.76%	1.148%
2	3.179	20	24	29	rVB2	32794	53536	3.93%	0.463%
3	3.455	67	71	76	rVB2	6381	9468	0.70%	0.082%
4	3.748	117	121	124	rBV3	2263	3396	0.25%	0.029%
5	3.878	139	143	155	rBV	69004	138972	10.21%	1.202%
6	4.207	193	199	205	rBV4	6687	10929	0.80%	0.094%
7	6.128	521	526	528	rBV4	2960	3111	0.23%	0.027%
8	7.215	703	711	722	rBV	151097	279377	20.52%	2.416%
9	7.385	731	740	748	rBV	92395	166662	12.24%	1.441%
10	7.591	768	775	785	rBV	135090	248997	18.29%	2.153%
11	8.061	847	855	861	rBV	81083	144420	10.61%	1.249%
12	8.120	861	865	870	rVV4	2758	5454	0.40%	0.047%
13	8.766	968	975	984	rBV2	190103	341550	25.09%	2.953%
14	9.230	1046	1054	1062	rBV	150252	273404	20.08%	2.364%
15	9.636	1117	1123	1127	rVB2	2365	3312	0.24%	0.029%
16	9.959	1170	1178	1188	rBV	138655	253929	18.65%	2.196%
17	10.499	1263	1270	1281	rBV	189282	354553	26.05%	3.066%
18	10.640	1289	1294	1305	rBV3	20311	38994	2.86%	0.337%
19	10.881	1328	1335	1343	rVB	109563	199498	14.66%	1.725%
20	11.016	1350	1358	1366	rBV	178867	332526	24.43%	2.875%
21	12.474	1604	1606	1610	rVB3	2116	2530	0.19%	0.022%
22	13.531	1782	1786	1790	rBV4	1554	2691	0.20%	0.023%
23	14.107	1878	1884	1891	rBV	359865	538501	39.56%	4.656%
24	14.407	1927	1935	1944	rVB	367312	599336	44.03%	5.182%
25	14.712	1980	1987	1994	rVB	170223	271394	19.94%	2.347%
26	14.906	2011	2020	2031	rBV2	144631	252728	18.57%	2.185%
27	15.705	2149	2156	2167	rBV	471125	737607	54.19%	6.378%
28	15.817	2167	2175	2185	rBV	147615	224761	16.51%	1.943%
29	15.940	2193	2196	2205	rVB4	5238	8971	0.66%	0.078%
30	17.321	2427	2431	2432	rBV3	3462	3847	0.28%	0.033%
31	17.462	2449	2455	2461	rBV	216315	340756	25.03%	2.946%
32	17.562	2464	2472	2480	rVB	544822	828007	60.83%	7.160%
33	18.337	2599	2604	2610	rBV5	13638	19307	1.42%	0.167%
34	18.772	2671	2678	2685	rBV	35937	58761	4.32%	0.508%
35	18.919	2698	2703	2707	rBV4	2476	3781	0.28%	0.033%
36	19.277	2760	2764	2767	rVB3	2101	3130	0.23%	0.027%

LSC Area Percent Report

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Sample : M2970-02
Misc :
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Instrument : BNA_G
ClientSampleId : BGE81

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_G\METHODS\SFAM-EPA-BG070921.M
Title : SVOA CALIBRATION

37	19.407	2782	2786	2791	rBV7	6586	10480	0.77%	0.091%
38	19.483	2794	2799	2804	rBV2	7430	11816	0.87%	0.102%
39	19.606	2818	2820	2824	rVB5	3421	4287	0.31%	0.037%
40	19.847	2851	2861	2870	rBV	669868	966386	70.99%	8.356%
41	20.241	2924	2928	2931	rVB5	2119	3232	0.24%	0.028%
42	20.347	2940	2946	2947	rBV4	5637	8572	0.63%	0.074%
43	20.593	2986	2988	2991	rVV3	3538	3902	0.29%	0.034%
44	20.746	3010	3014	3017	rBV4	4793	5981	0.44%	0.052%
45	20.817	3021	3026	3037	rVV	1161308	1361241	100.00%	11.770%
46	20.928	3042	3045	3046	rBV2	2878	2868	0.21%	0.025%
47	21.011	3057	3059	3062	rVB4	5158	4727	0.35%	0.041%
48	21.316	3109	3111	3112	rBV2	3733	2627	0.19%	0.023%
49	21.369	3115	3120	3126	rBV2	53976	92098	6.77%	0.796%
50	21.557	3150	3152	3156	rVB5	4260	3537	0.26%	0.031%
51	21.616	3160	3162	3166	rVB4	3252	3096	0.23%	0.027%
52	21.739	3177	3183	3191	rVB	242392	389003	28.58%	3.364%
53	21.927	3213	3215	3216	rBV	4572	3233	0.24%	0.028%
54	22.027	3229	3232	3233	rVB3	4418	3527	0.26%	0.030%
55	22.503	3310	3313	3314	rBV2	2813	3217	0.24%	0.028%
56	22.568	3317	3324	3329	rBV3	30103	68682	5.05%	0.594%
57	22.855	3372	3373	3377	rVB3	3912	2969	0.22%	0.026%
58	22.920	3382	3384	3387	rBV3	3402	3309	0.24%	0.029%
59	23.096	3412	3414	3416	rBV3	3490	3176	0.23%	0.027%
60	23.261	3437	3442	3447	rVB8	8408	16831	1.24%	0.146%
61	23.302	3447	3449	3452	rVB4	3268	2868	0.21%	0.025%
62	23.420	3466	3469	3470	rBV3	3087	2670	0.20%	0.023%
63	23.455	3470	3475	3479	rVV6	7626	13420	0.99%	0.116%
64	23.737	3521	3523	3526	rVB4	3192	2849	0.21%	0.025%
65	23.766	3526	3528	3530	rBV3	3287	2953	0.22%	0.026%
66	23.931	3551	3556	3561	rBV8	10857	19612	1.44%	0.170%
67	24.083	3576	3582	3589	rVB4	16965	35970	2.64%	0.311%
68	24.154	3589	3594	3596	rBV5	7536	14192	1.04%	0.123%
69	24.365	3627	3630	3632	rBV4	2545	3044	0.22%	0.026%
70	24.659	3676	3680	3682	rVB5	3255	4061	0.30%	0.035%
71	24.806	3691	3705	3715	rBV2	362210	984276	72.31%	8.511%
72	25.029	3732	3743	3757	rVB2	145583	446138	32.77%	3.858%
73	25.241	3776	3779	3781	rBV4	2519	2533	0.19%	0.022%
74	25.329	3793	3794	3799	rBV5	3389	4006	0.29%	0.035%
75	25.658	3848	3850	3853	rBV3	2109	2647	0.19%	0.023%
76	25.699	3855	3857	3860	rVB4	3759	3214	0.24%	0.028%

LSC Area Percent Report

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Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

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
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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Title : SVOA CALIBRATION

77	26.022	3909	3912	3913	rBV3	2631	2962	0.22%	0.026%
78	26.216	3936	3945	3956	rVB5	21295	67575	4.96%	0.584%
79	26.340	3964	3966	3967	rBV	4308	2955	0.22%	0.026%
80	26.557	4000	4003	4005	rVB3	3254	3930	0.29%	0.034%
81	27.321	4132	4133	4138	rVB5	3399	2911	0.21%	0.025%
82	27.456	4155	4156	4162	rBV5	3807	4305	0.32%	0.037%
83	28.114	4266	4268	4272	rBV5	2087	2573	0.19%	0.022%
84	28.232	4286	4288	4291	rBV4	2437	3209	0.24%	0.028%
85	28.355	4305	4309	4312	rBV6	2825	3630	0.27%	0.031%
86	28.625	4351	4355	4356	rBV4	3826	3525	0.26%	0.030%
87	28.931	4404	4407	4409	rBV4	2520	3301	0.24%	0.029%
88	29.271	4463	4465	4467	rBV2	6862	7640	0.56%	0.066%
89	29.695	4535	4537	4539	rBV3	2883	3056	0.22%	0.026%
90	29.847	4562	4563	4566	rBV3	3655	3472	0.26%	0.030%
91	30.282	4635	4637	4639	rVV3	3437	2781	0.20%	0.024%
92	30.317	4641	4643	4646	rVB4	3358	3096	0.23%	0.027%
93	30.417	4658	4660	4665	rVB5	2984	4552	0.33%	0.039%
94	30.535	4677	4680	4682	rVB3	5344	4542	0.33%	0.039%
95	30.828	4726	4730	4731	rVB3	2752	2604	0.19%	0.023%
96	30.846	4731	4733	4735	rVB3	4495	2773	0.20%	0.024%
97	31.487	4840	4842	4846	rBV4	3281	4949	0.36%	0.043%
98	31.816	4896	4898	4901	rBV4	2602	2839	0.21%	0.025%
99	32.280	4975	4977	4980	rBV3	2206	2931	0.22%	0.025%
100	32.620	5031	5035	5038	rBV4	3434	4727	0.35%	0.041%

Sum of corrected areas: 11565082

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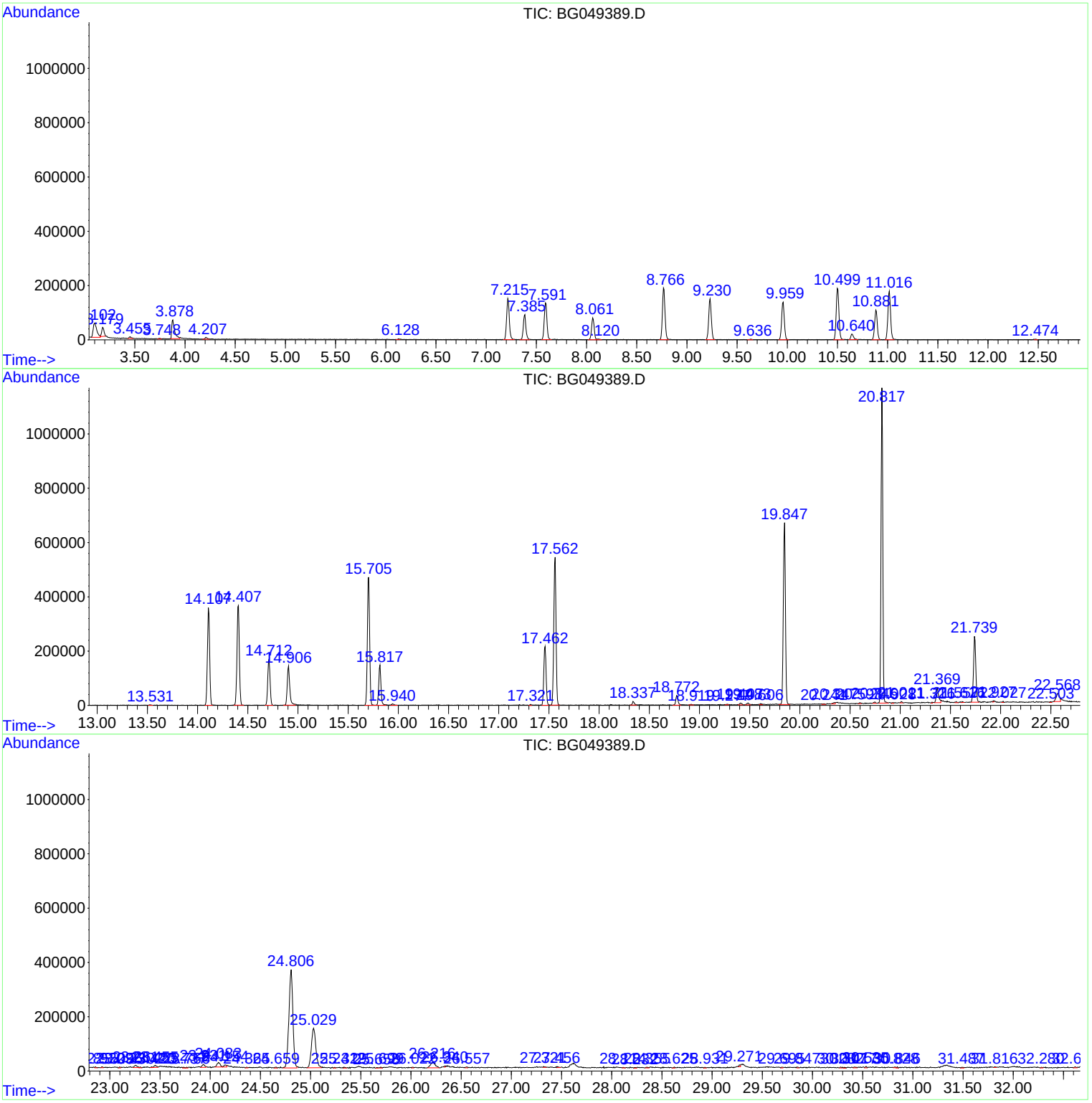
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGE81

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071621\
Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
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Sample : M2970-02
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Instrument : BNA_G
ClientSampleId : BGE81

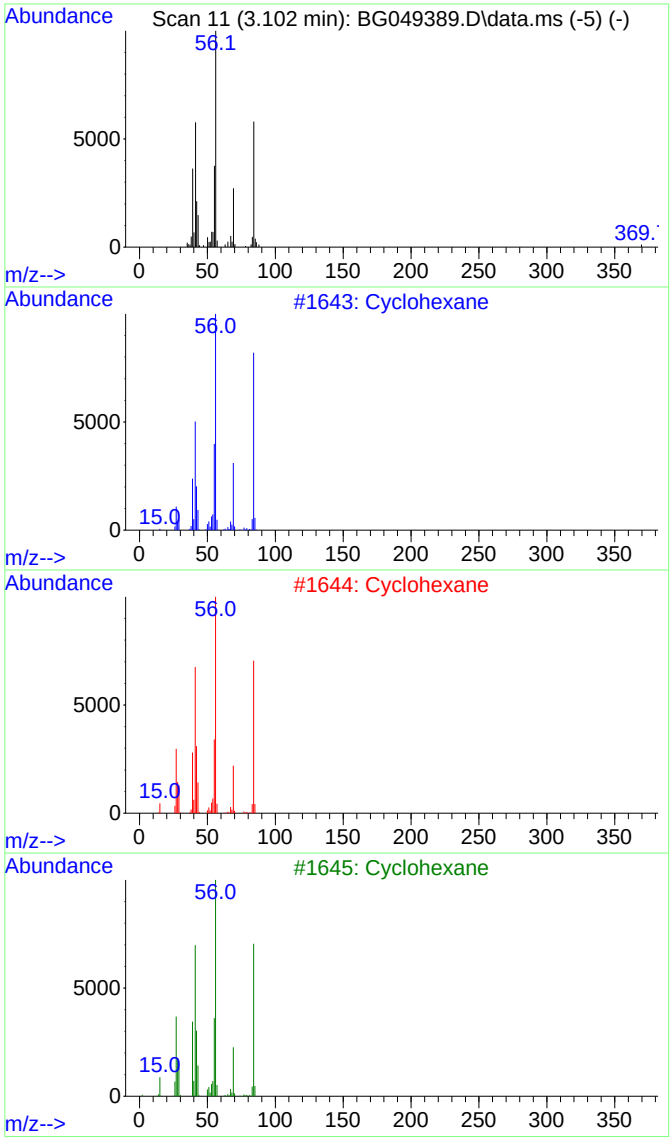
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

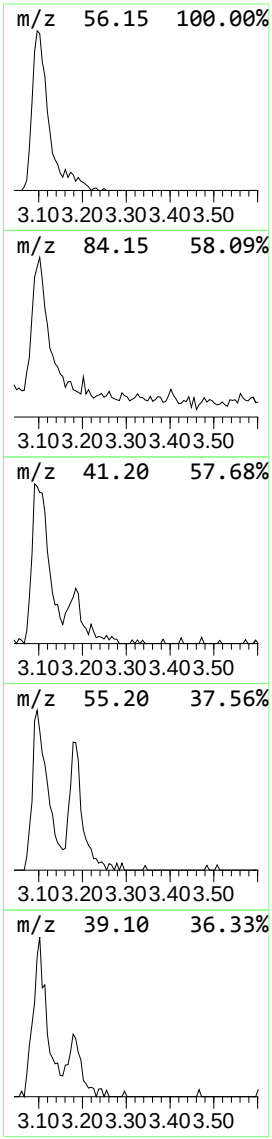
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.10 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.100	18.39 ng/ul	132798	1,4-Dichlorobenzene-d4	8.067

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	87
2	Cyclohexane	84	C6H12	000110-82-7	87
3	Cyclohexane	84	C6H12	000110-82-7	87
4	Cyclohexane	84	C6H12	000110-82-7	86
5	Cyclohexane	84	C6H12	000110-82-7	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071621\
Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

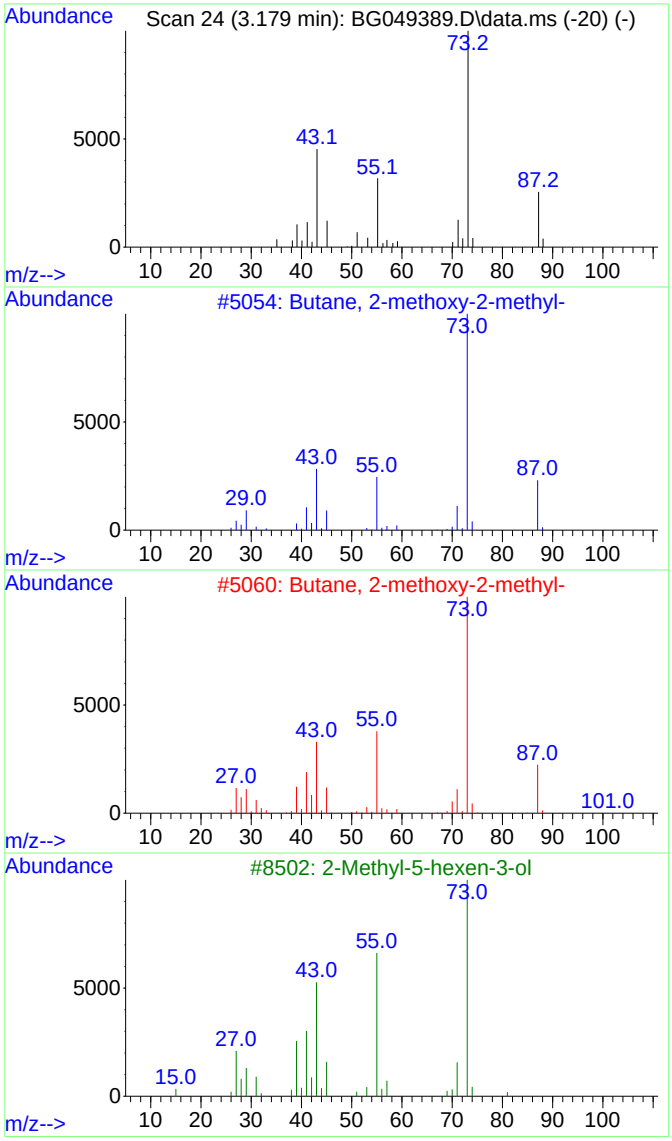
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

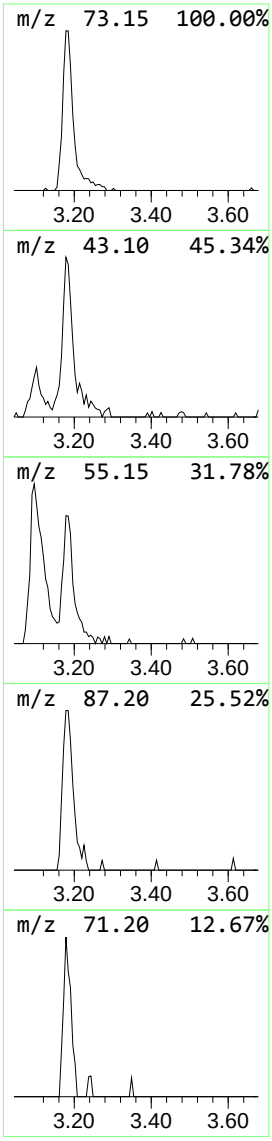
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.180	7.41 ng/ul	53536	1,4-Dichlorobenzene-d4	8.067

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23
4	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9
5	1,4-Butanediamine	88	C4H12N2	000110-60-1	9





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Instrument : BNA_G
ClientSampleId : BGE81

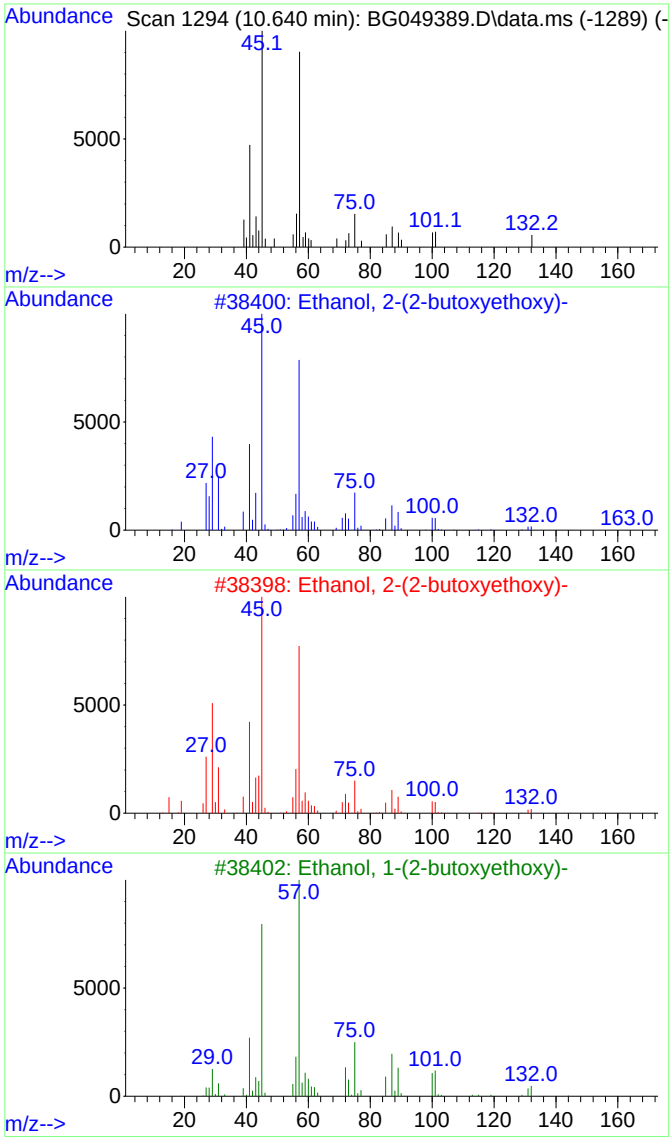
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

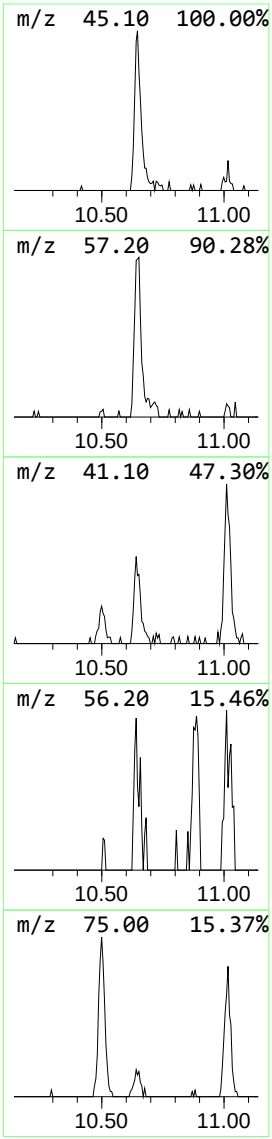
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.640	3.91 ng/ul	38994	Naphthalene-d8	10.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	64
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
5	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	64





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

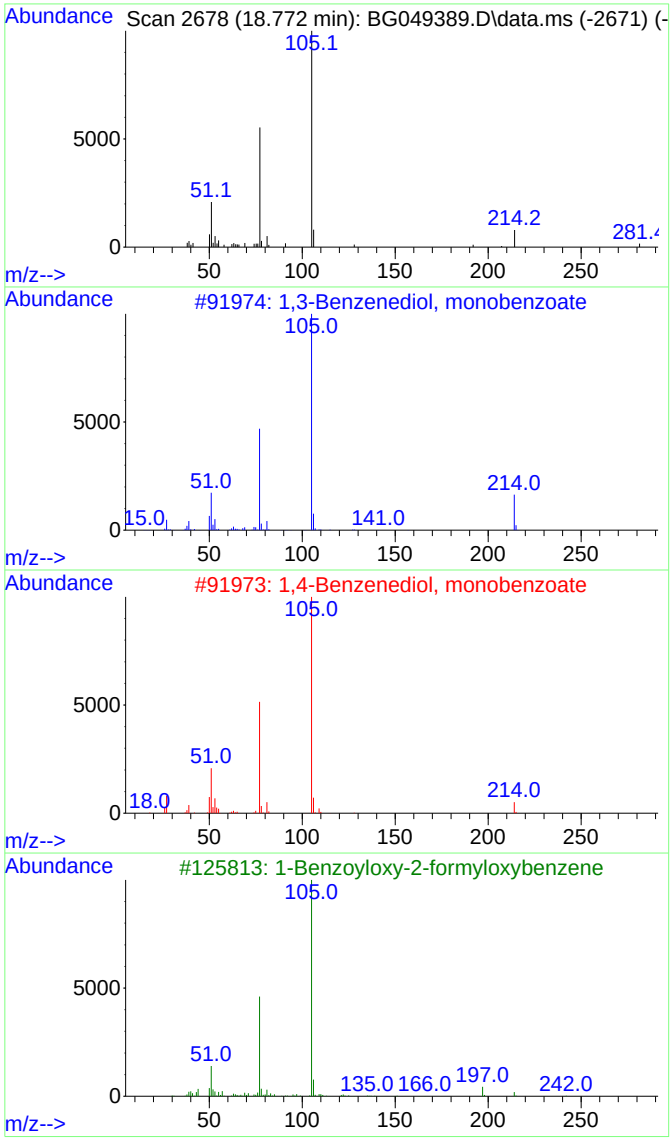
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

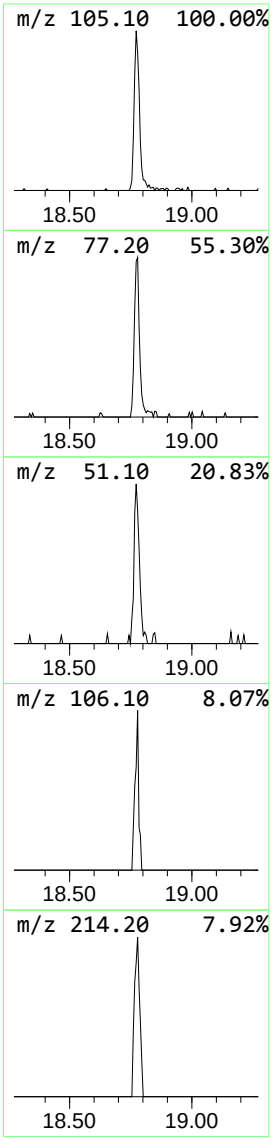
****

Peak Number 4 1,3-Benzenediol, monobenzoate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.770	3.45 ng/ul	58761	Phenanthrene-d10	17.462

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	87
2	1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	86
3	1-Benzoyloxy-2-formyloxybenzene	242	C14H10O4	079792-93-1	78
4	1,2-benzenediol, monobenzoate	214	C13H10O3	1000400-67-6	78
5	N-Methylbenzamide, N-pentafluoro...	281	C11H8F5NO2	1000446-98-1	72





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Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

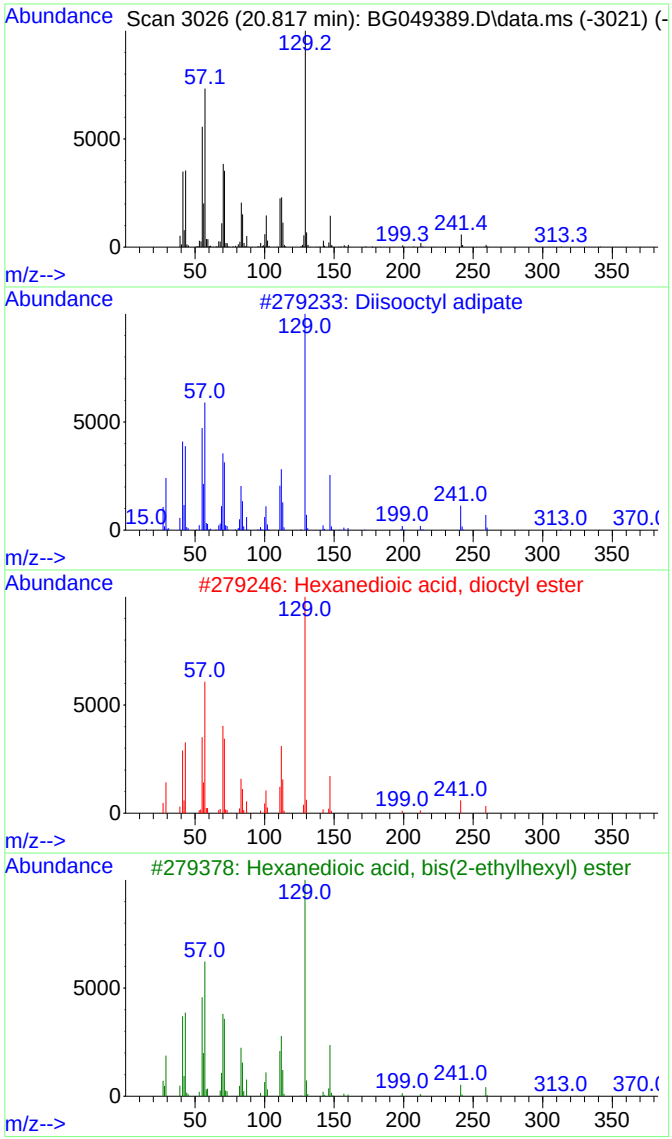
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

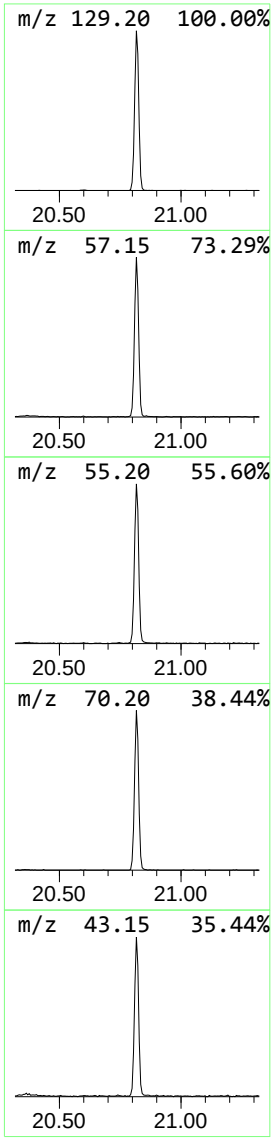
****

Peak Number 5 Diisooctyl adipate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.820	69.99 ng/ul	1361240	Chrysene-d12	21.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisooctyl adipate	370	C22H42O4	001330-86-5	91
2	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	90
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	74
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	72
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	72





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Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

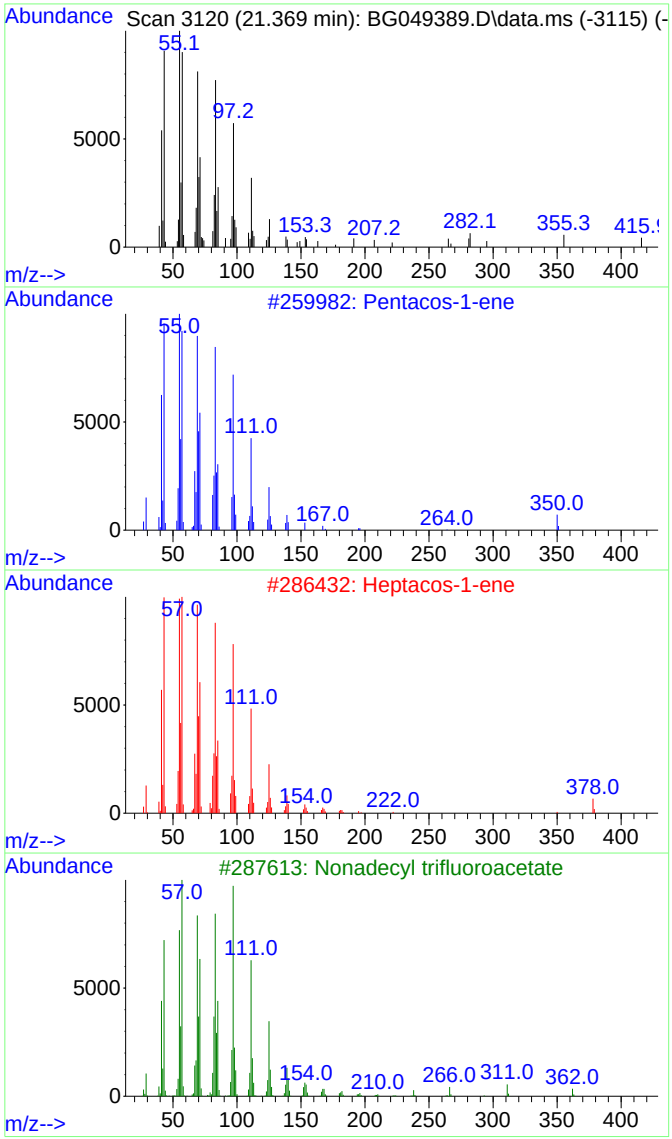
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

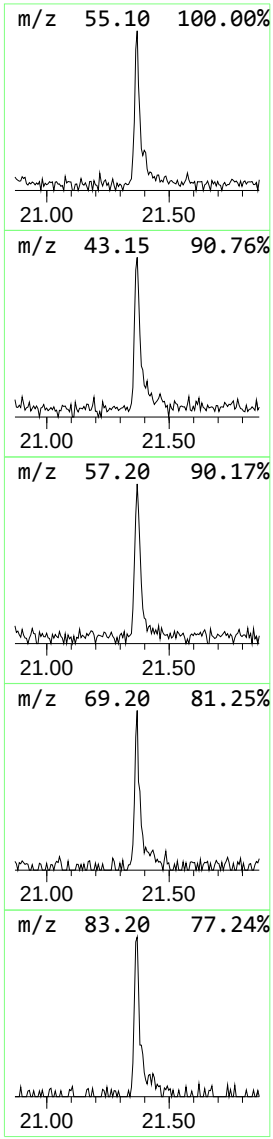
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.370	4.74 ng/ul	92098	Chrysene-d12	21.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacos-1-ene	350	C25H50	016980-85-1	90
2	Heptacos-1-ene	378	C27H54	015306-27-1	87
3	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	74
4	1-Tricosene	322	C23H46	018835-32-0	58
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	50





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Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

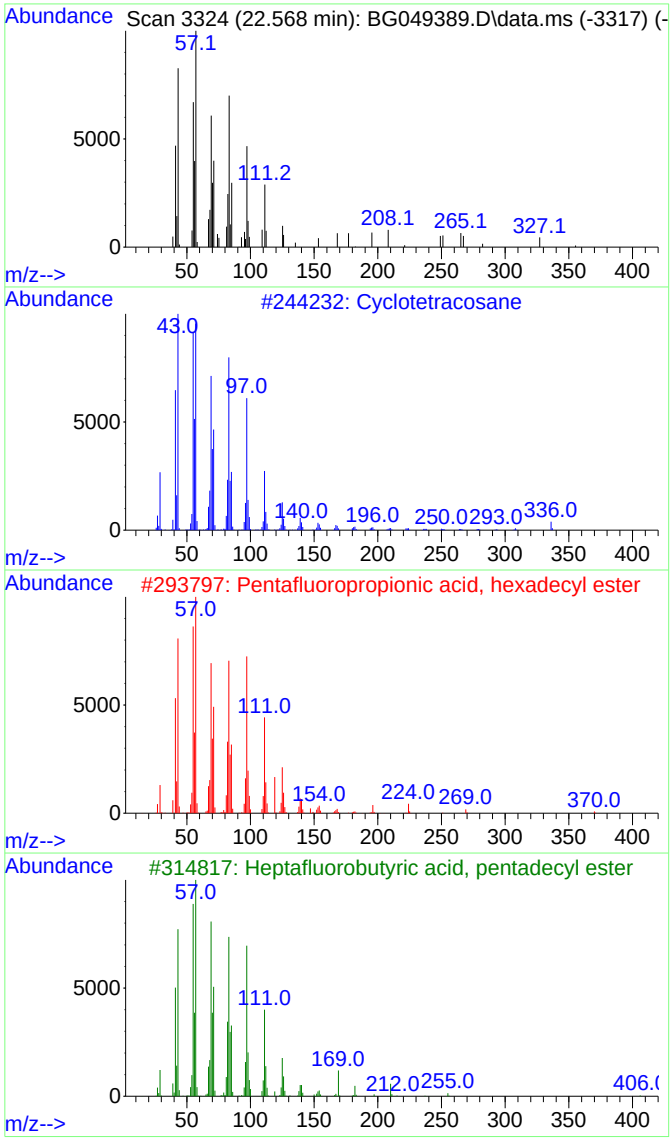
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

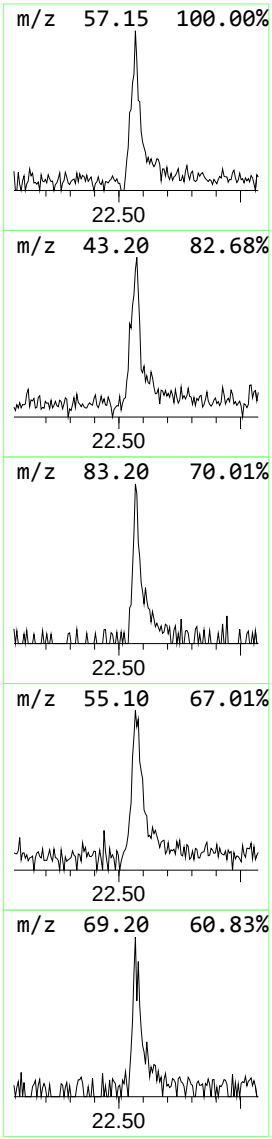
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic22.57 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.570	3.53 ng/ul	68682	Chrysene-d12	21.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	90
2	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	87
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	87
4	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87





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Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

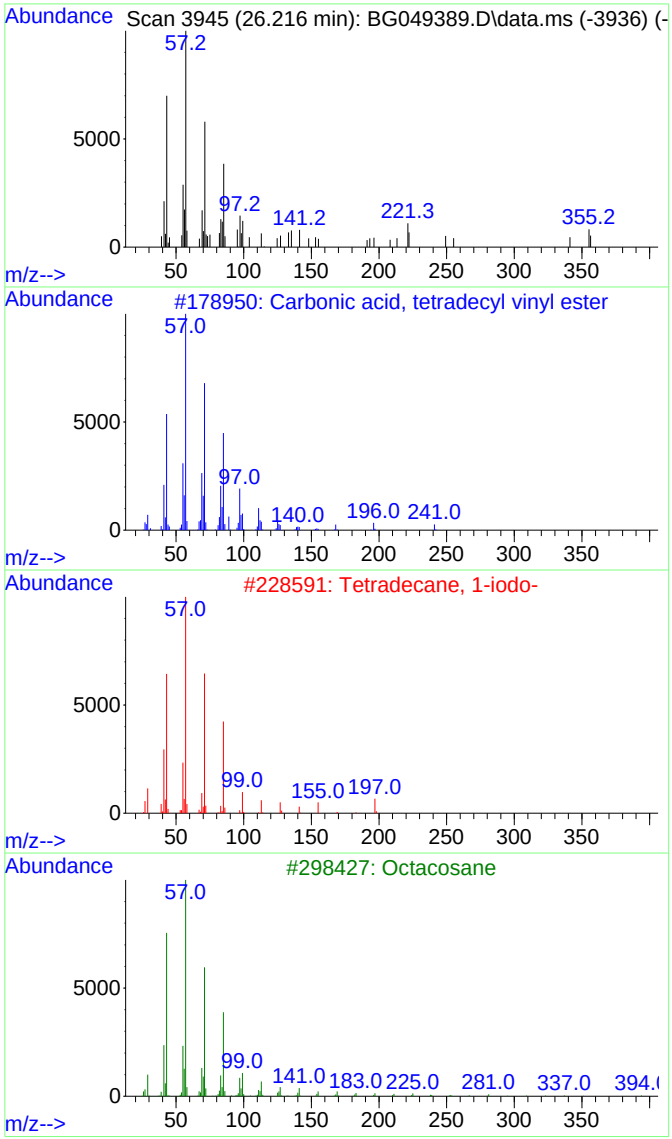
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

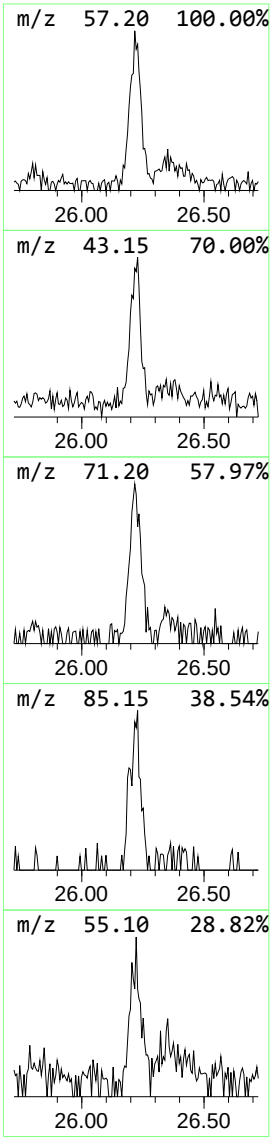
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Carbonic acid, tetradecyl v... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.220	3.03 ng/ul	67575	Perylene-d12	25.029

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	52
2	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	49
3	Octacosane	394	C28H58	000630-02-4	49
4	Pentadecane	212	C15H32	000629-62-9	49
5	Octadecane	254	C18H38	000593-45-3	49





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071621\
Data File : BG049389.D
Acq On : 16 Jul 2021 22:18
Operator : CG/JU
Sample : M2970-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGE81

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.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
(DEL) Alkane: C...	3.100	18.4	ng/ul	132798	1	8.067	144420	20.0
Butane, 2-metho...	3.180	7.4	ng/ul	53536	1	8.067	144420	20.0
Ethanol, 2-(2-b...	10.640	3.9	ng/ul	38994	2	10.881	199498	20.0
1,3-Benzenediol...	18.770	3.5	ng/ul	58761	4	17.462	340756	20.0
Diisooctyl adipate	20.820	70.0	ng/ul	1361240	5	21.739	389003	20.0
(DEL) Alkane: S...	21.370	4.7	ng/ul	92098	5	21.739	389003	20.0
(DEL) Alkane: C...	22.570	3.5	ng/ul	68682	5	21.739	389003	20.0
Carbonic acid, ...	26.220	3.0	ng/ul	67575	6	25.029	446138	20.0