LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

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-BG080421.M
Title : SVOA CALIBRATION

Signal : TIC: BG049647.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.074	3	6	15	rVB3	64165	120373	3.72%	0.942%
2	3.156	15	20	31	rVB	52854	95509	2.95%	0.747%
3	3.426	62	66	73	rVB6	6710	12699	0.39%	0.099%
4	3.849	134	138	155	rBV	78424	166049	5.13%	1.299%
5	7.192	699	707	720	rVB	146925	281122	8.68%	2.199%
6	7.362	730	736	747	rVB	93811	162617	5.02%	1.272%
7	7.568	765	771	783	rVB	141464	254544	7.86%	1.991%
8	8.044	845	852	857	rBV2	79541	140006	4.32%	1.095%
9	8.096	857	861	866	rVB3	9076	14513	0.45%	0.114%
10	8.749	964	972	980	rBV	181418	325670	10.06%	2.548%
11	9.213	1041	1051	1058	rBV	148195	268067	8.28%	2.097%
12	9.612	1116	1119	1125	rVB3	2686	4308	0.13%	0.034%
13	9.935	1166	1174	1188	rBV2	130102	244847	7.56%	1.916%
14	10.153	1203	1211	1213	rBV4	4282	8606	0.27%	0.067%
15	10.481	1260	1267	1278	rBV	181871	333464	10.30%	2.609%
16	10.628	1284	1292	1304	rBV3	21575	46992	1.45%	0.368%
17	10.863	1324	1332	1340	rVB	111375	198659	6.14%	1.554%
18	10.998	1347	1355	1363	rBV2	159439	301472	9.31%	2.359%
19	14.094	1875	1882	1888	rBV	331478	494517	15.27%	3.869%
20	14.388	1925	1932	1940	rBV	348193	551813	17.04%	4.317%
21	14.693	1976	1984	1991	rBV	178604	274396	8.47%	2.147%
22	14.887	2011	2017	2032	rBV	135362	247754	7.65%	1.938%
23	15.686	2146	2153	2159	rBV	443921	667046	20.60%	5.219%
24	15.804	2167	2173	2180	rBV2	162643	244381	7.55%	1.912%
25	15.927	2190	2194	2199	rVB4	5346	8248	0.25%	0.065%
26	17.443	2445	2452	2459	rBV	221949	333462	10.30%	2.609%
27	17.543	2462	2469	2476	rBV2	443129	704791	21.77%	5.514%
28	18.101	2559	2564	2568	rBV6	5753	9161	0.28%	0.072%
29	18.153	2568	2573	2577	rVB4	5787	7617	0.24%	0.060%
30	18.324	2598	2602	2609	rBV3	20880	33059	1.02%	0.259%
31	18.759	2670	2676	2684	rBV	126764	187968	5.81%	1.471%
32	19.240	2754	2758	2761	rBV4	3533	5021	0.16%	0.039%
33	19.393	2780	2784	2787	rBV3	7022	8789	0.27%	0.069%
34	19.475	2791	2798	2801	rBV4	7377	14440	0.45%	0.113%
35	19.593	2813	2818	2822	rBV4	7960	11629	0.36%	0.091%
36	19.834	2850	2859	2865	rBV	544685	793521	24.51%	6.208%

LSC Area Percent Report

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Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

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-BG080421.M
Title : SVOA CALIBRATION

37	20.345	2944	2946	2953	rVB5	5908	8374	0.26%	0.066%
38	20.580	2984	2986	2990	rVB4	7900	8396	0.26%	0.066%
39	20.803	3019	3024	3029	rBV	2603797	3238026	100.00%	25.333%
40	20.844	3030	3031	3035	rVB3	5696	4393	0.14%	0.034%
41	21.032	3060	3063	3064	rBV3	3062	3243	0.10%	0.025%
42	21.173	3084	3087	3089	rVB3	4142	4329	0.13%	0.034%
43	21.355	3113	3118	3124	rBV3	41986	75578	2.33%	0.591%
44	21.725	3175	3181	3188	rBV	213360	341931	10.56%	2.675%
45	21.907	3210	3212	3216	rVB3	5775	7600	0.23%	0.059%
46	22.530	3313	3318	3324	rBV5	16911	39130	1.21%	0.306%
47	23.211	3428	3434	3442	rVB7	10128	28911	0.89%	0.226%
48	23.335	3453	3455	3459	rBV4	2944	3446	0.11%	0.027%
49	23.429	3465	3471	3476	rBV9	6476	11272	0.35%	0.088%
50	23.646	3506	3508	3513	rVB5	2695	3392	0.10%	0.027%
51	23.904	3545	3552	3560	rVB5	40096	90856	2.81%	0.711%
52	24.051	3570	3577	3586	rVB5	17788	43386	1.34%	0.339%
53	24.134	3586	3591	3592	rBV5	7425	12340	0.38%	0.097%
54	24.774	3687	3700	3713	rVB2	272796	762463	23.55%	5.965%
55	25.003	3728	3739	3752	rBV2	129006	373234	11.53%	2.920%
56	25.444	3807	3814	3821	rBV9	15819	42717	1.32%	0.334%
57	25.761	3861	3868	3872	rBV9	6369	12659	0.39%	0.099%
58	25.837	3878	3881	3883	rBV3	2627	3325	0.10%	0.026%
59	26.172	3931	3938	3943	rBV7	13267	33908	1.05%	0.265%
60	27.241	4118	4120	4124	rVV4	3249	6003	0.19%	0.047%
61	27.547	4169	4172	4174	rBV4	6238	6165	0.19%	0.048%
62	27.846	4219	4223	4226	rVB5	2983	4273	0.13%	0.033%
63	28.945	4409	4410	4417	rBV6	2373	3703	0.11%	0.029%
64	29.186	4449	4451	4452	rBV2	3712	3262	0.10%	0.026%
65	29.221	4453	4457	4458	rVV4	3139	4991	0.15%	0.039%
66	30.026	4592	4594	4598	rVB6	3588	4117	0.13%	0.032%
67	30.684	4702	4706	4707	rBV4	4120	3903	0.12%	0.031%
68	30.701	4707	4709	4712	rVB3	3945	3984	0.12%	0.031%
69	30.736	4712	4715	4717	rBV4	3603	4114	0.13%	0.032%
70	31.582	4857	4859	4862	rBV4	3474	3590	0.11%	0.028%
71	31.782	4889	4893	4894	rBV3	3049	3577	0.11%	0.028%

Sum of corrected areas: 12781721

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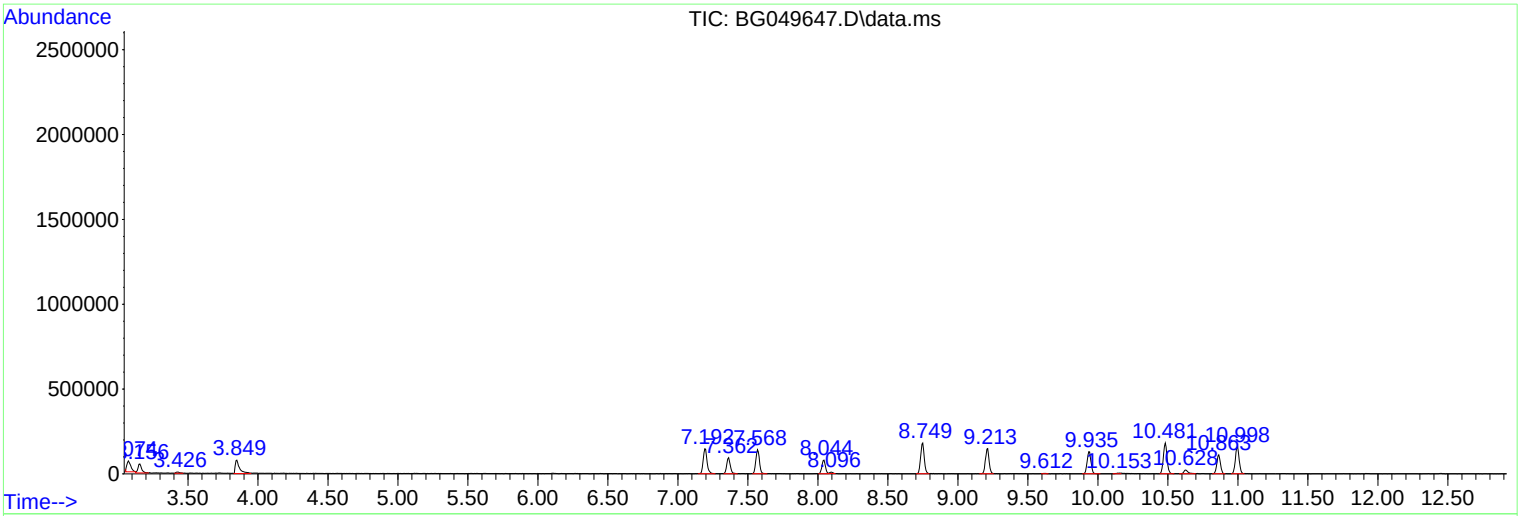
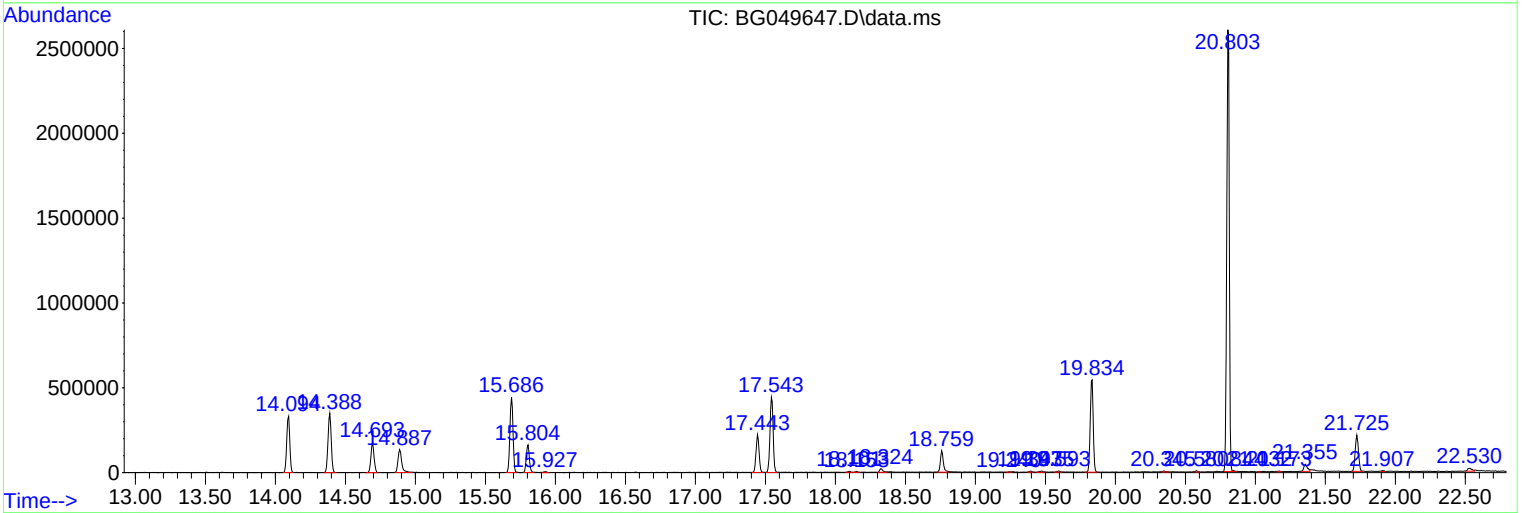
Misc :
ALS Vial : 8 Sample Multiplier: 1

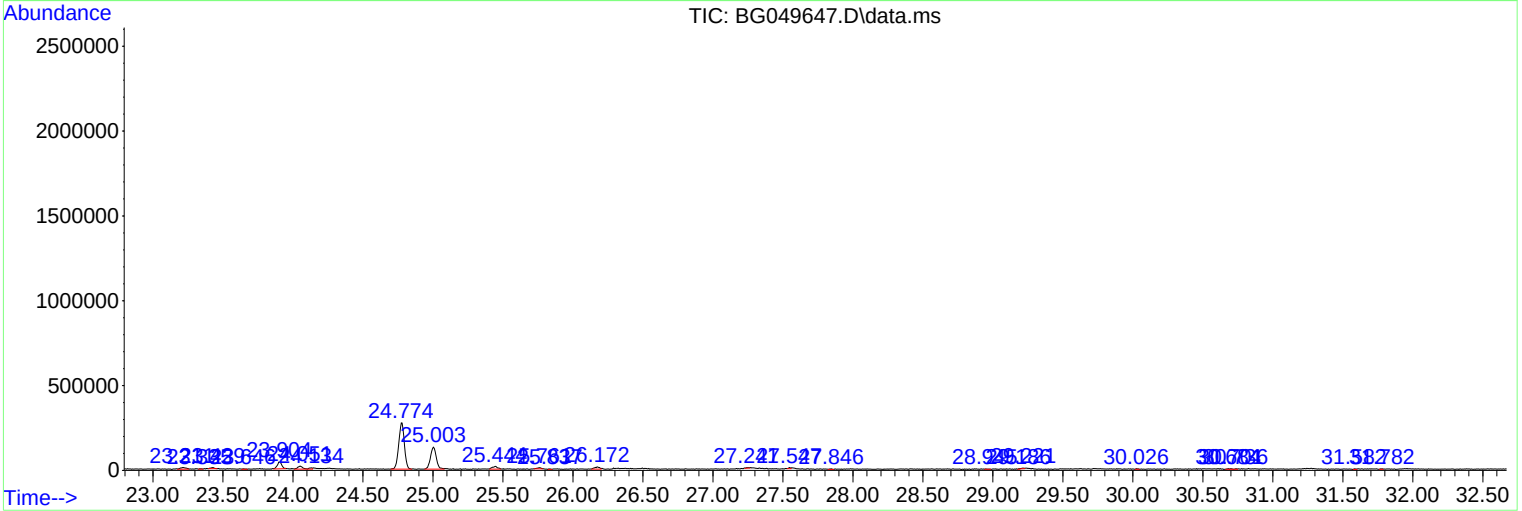
Instrument :
BNA_G
ClientSampleId :
BGET3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET3

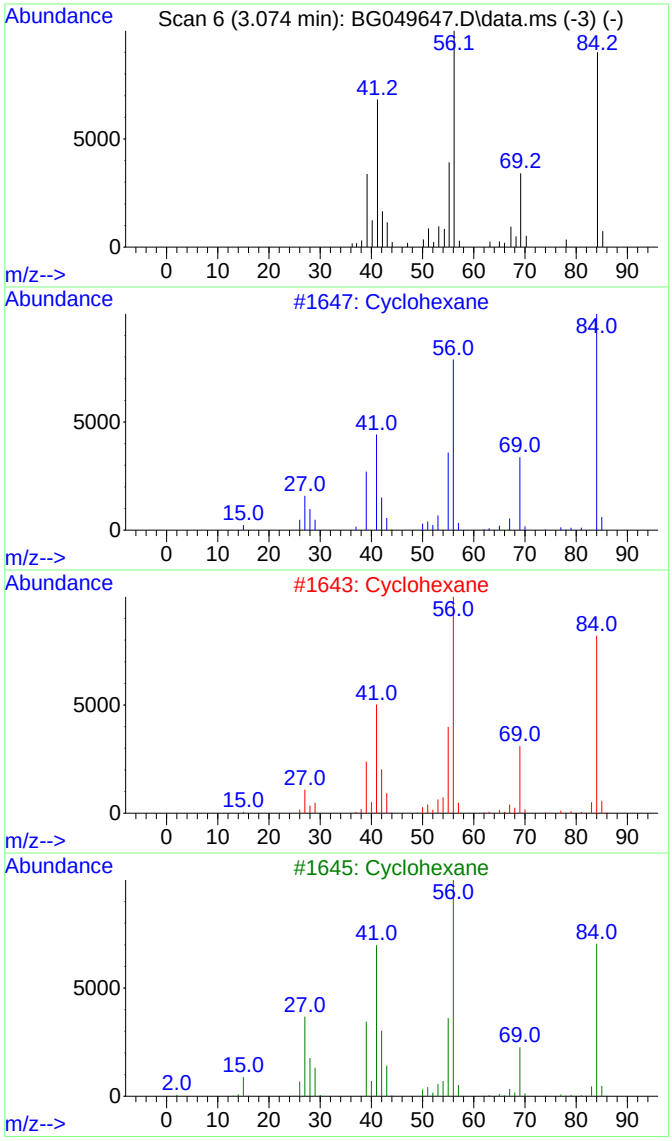
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

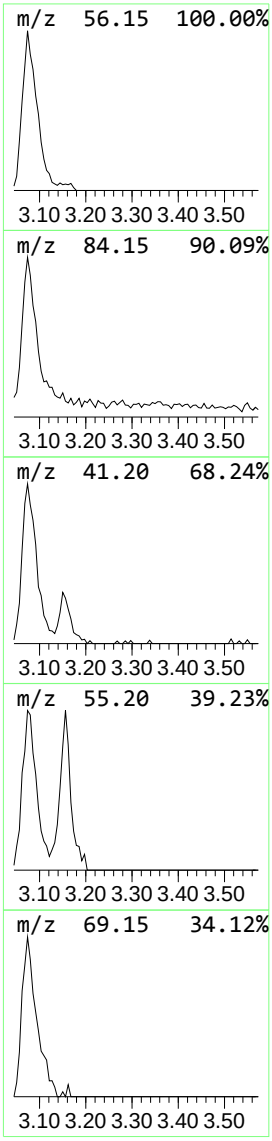
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.074 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.074	17.20 ng/ul	120373	1,4-Dichlorobenzene-d4	8.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	91
2	Cyclohexane	84	C6H12	000110-82-7	83
3	Cyclohexane	84	C6H12	000110-82-7	80
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	74
5	2-Amino-oxazole	84	C3H4N2O	004570-45-0	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

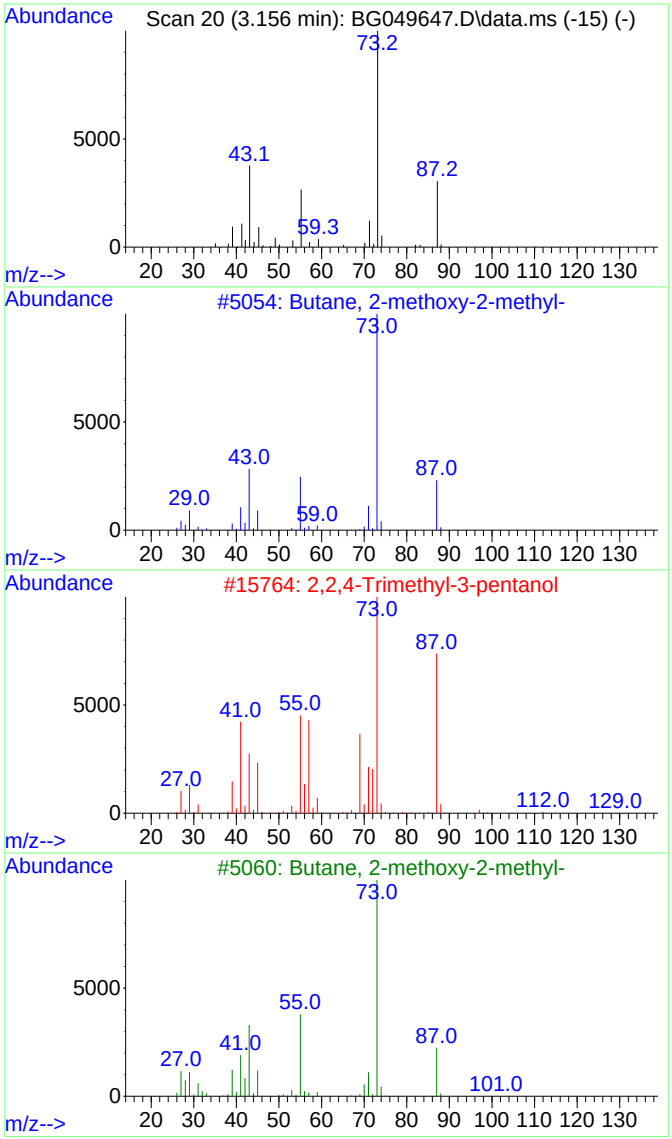
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

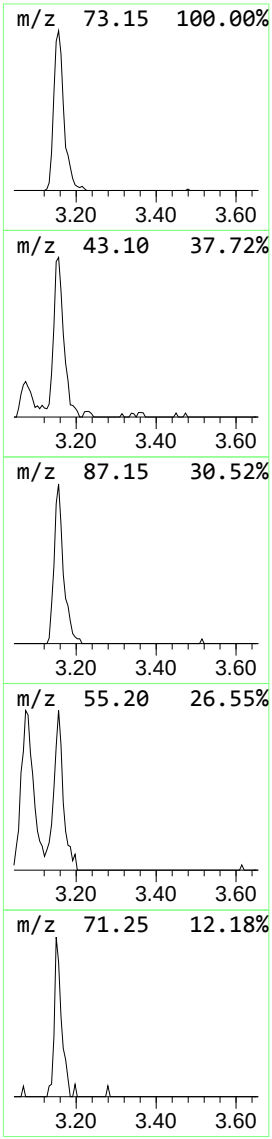
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.156	13.64 ng/ul	95509	1,4-Dichlorobenzene-d4	8.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	38
4	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	28
5	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25





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Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET3

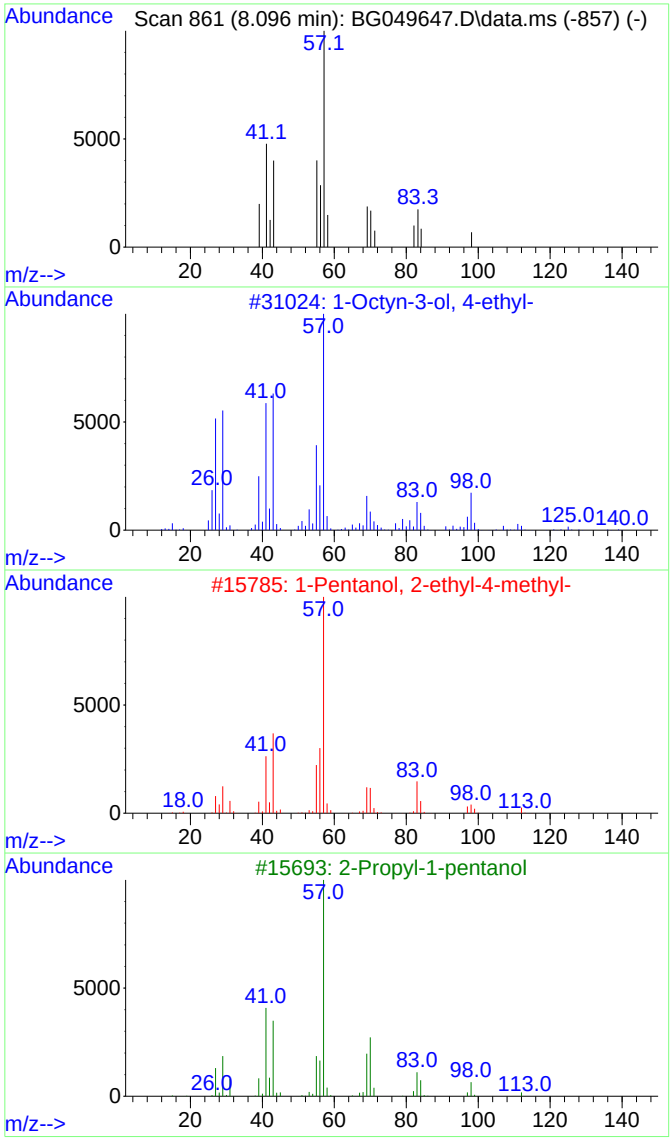
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

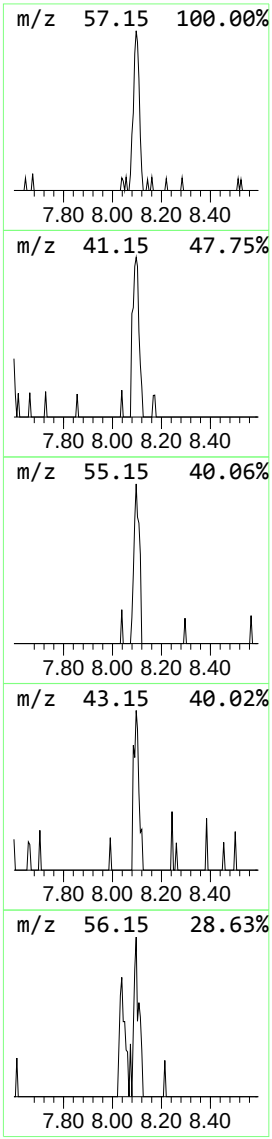
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Octyn-3-ol, 4-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.096	2.07 ng/ul	14513	1,4-Dichlorobenzene-d4	8.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octyn-3-ol, 4-ethyl-	154	C10H18O	005877-42-9	50
2	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	50
3	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	47
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	42
5	Hexane, 1-(hexyloxy)-4-methyl-	200	C13H28O	074421-20-8	38





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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

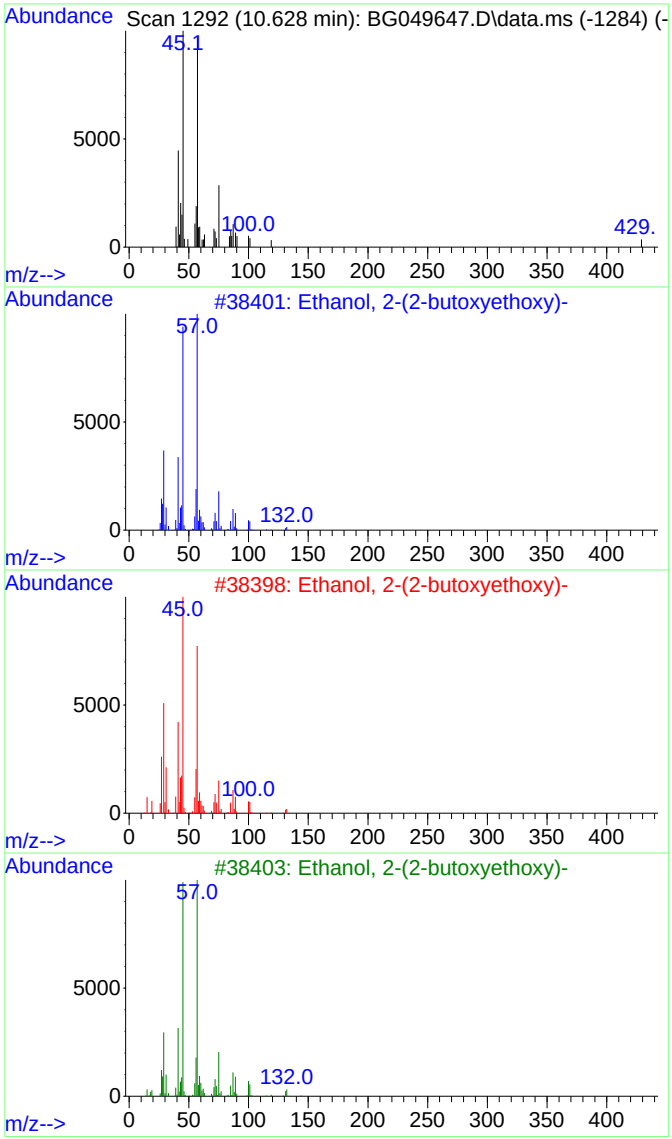
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

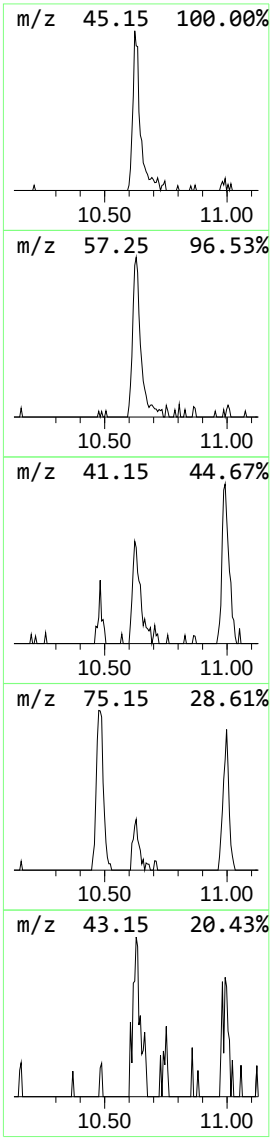
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.628	4.73 ng/ul	46992	Naphthalene-d8	10.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	72
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





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Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

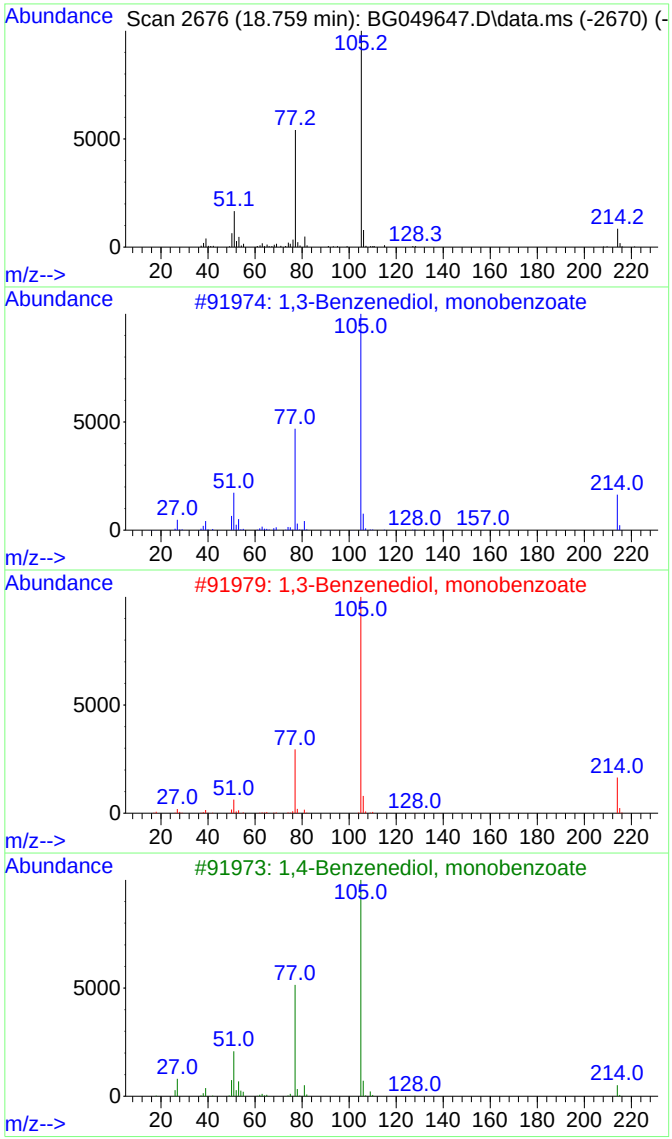
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

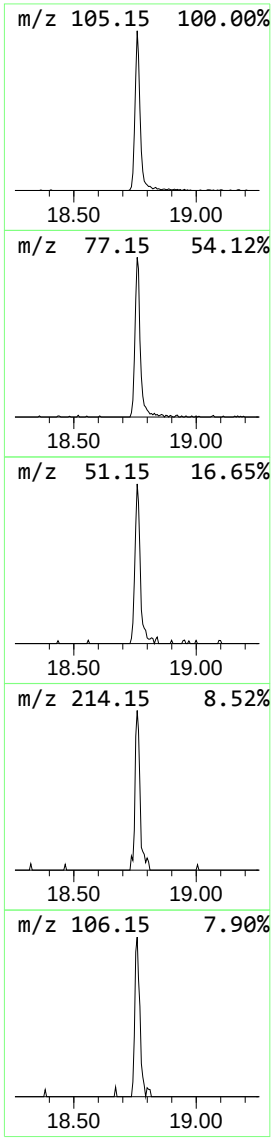
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,3-Benzenediol, monobenzoate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.759	11.27 ng/ul	187968	Phenanthrene-d10	17.443

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	87
2	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	86
3	1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	83
4	Benzamide, N-(1,2-dihydro-2-oxo-...	215	C11H9N3O2	026661-13-2	78
5	5-Oxo-5-phenylpentanenitrile	173	C11H11NO	1000486-83-7	72





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Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

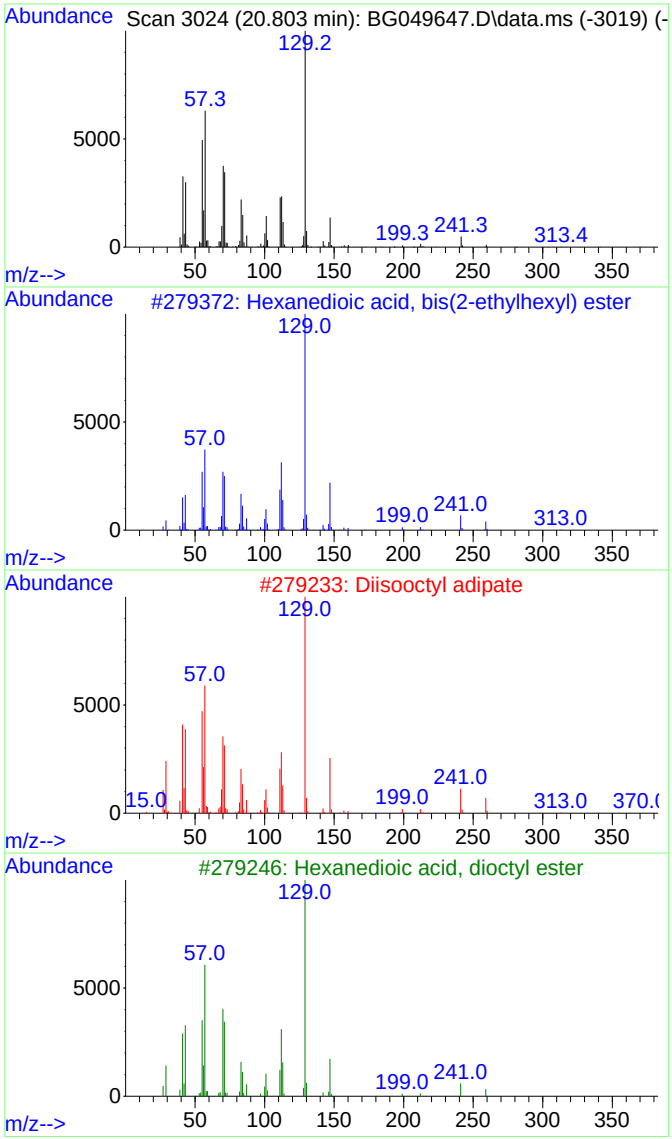
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

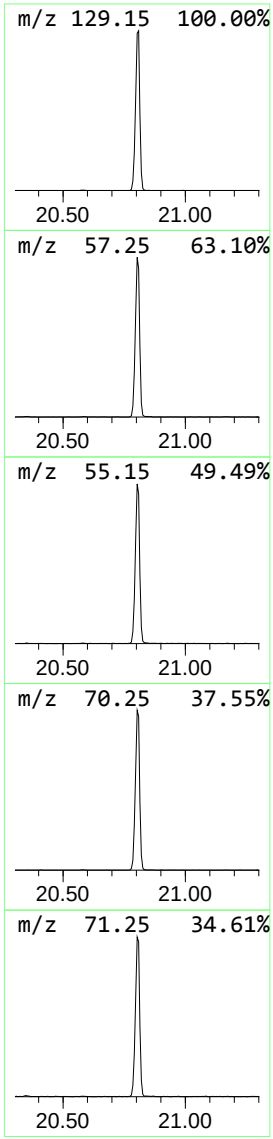
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.803	189.40 ng/ul	3238030	Chrysene-d12	21.725

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2	Diisooctyl adipate	370	C22H42O4	001330-86-5	91
3	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	90
4	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

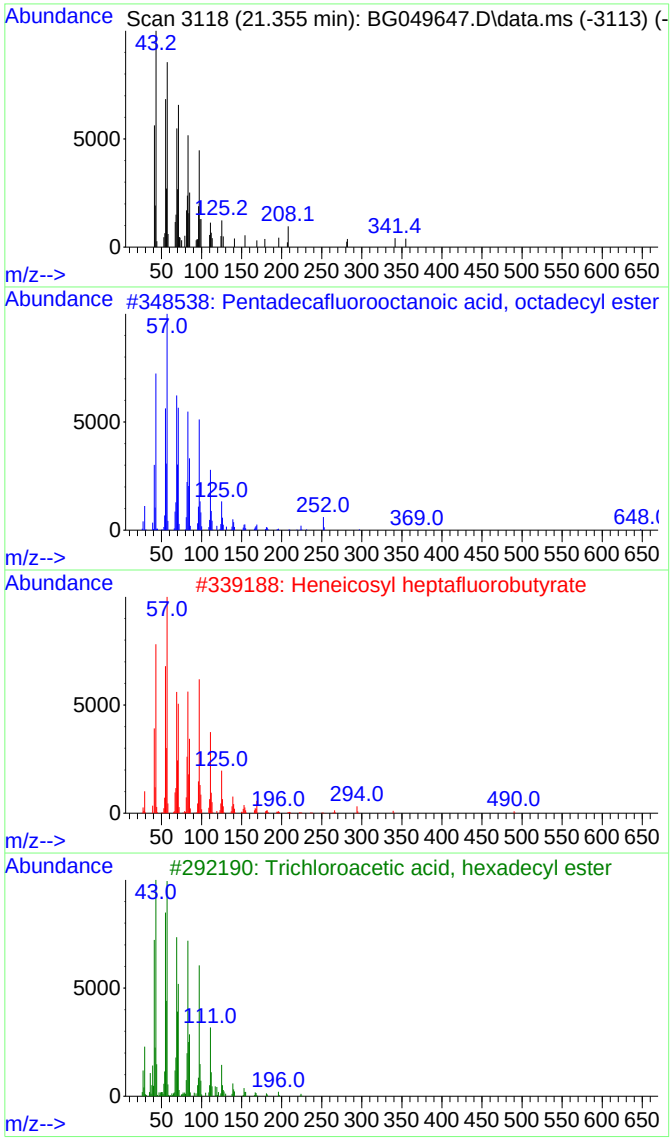
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

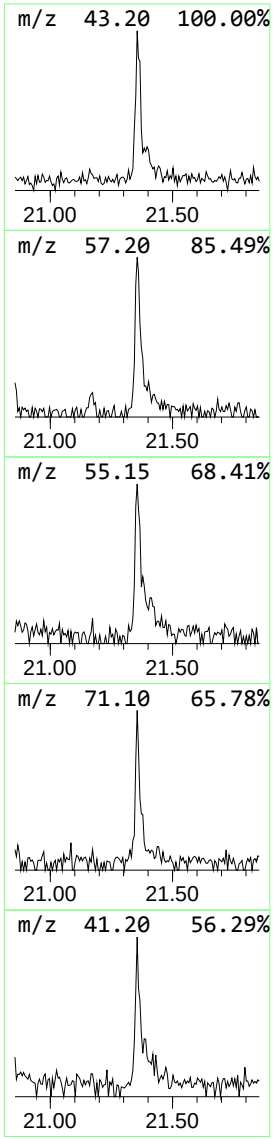
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.355	4.42 ng/ul	75578	Chrysene-d12	21.725

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	81
2	Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	74
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	62
4	17-Pentatriacontene	491	C35H70	006971-40-0	62
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	62





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Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

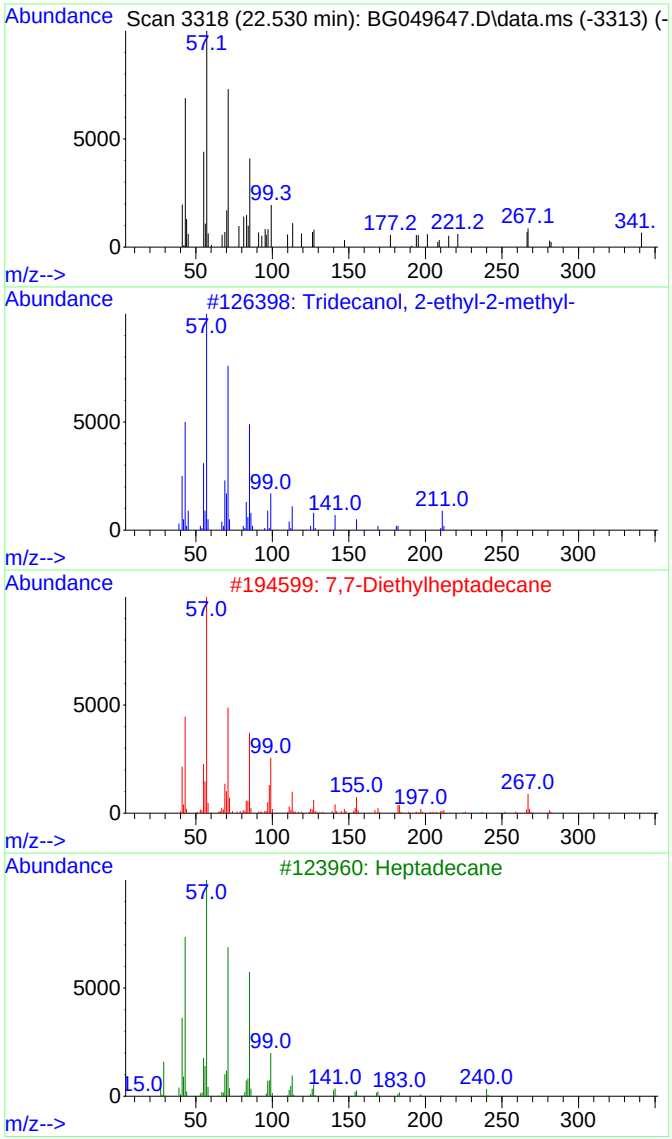
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

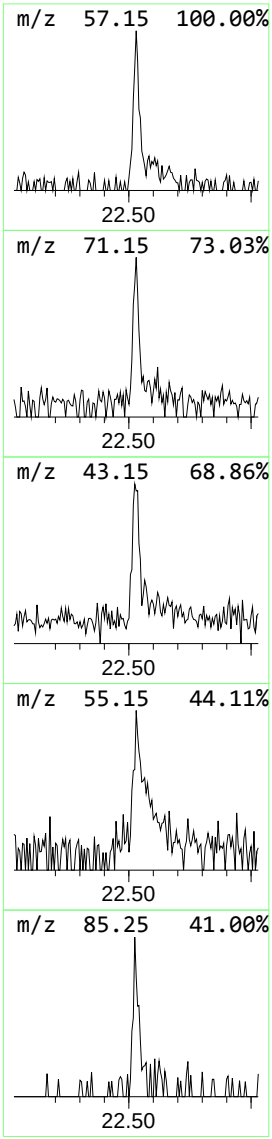
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Tridecanol, 2-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.530	2.29 ng/ul	39130	Chrysene-d12	21.725

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1010115-66-1	80
2	7,7-Diethylheptadecane	296	C21H44	1000360-41-5	72
3	Heptadecane	240	C17H36	000629-78-7	72
4	Triacontane	422	C30H62	000638-68-6	72
5	Octacosane	394	C28H58	000630-02-4	72





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Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

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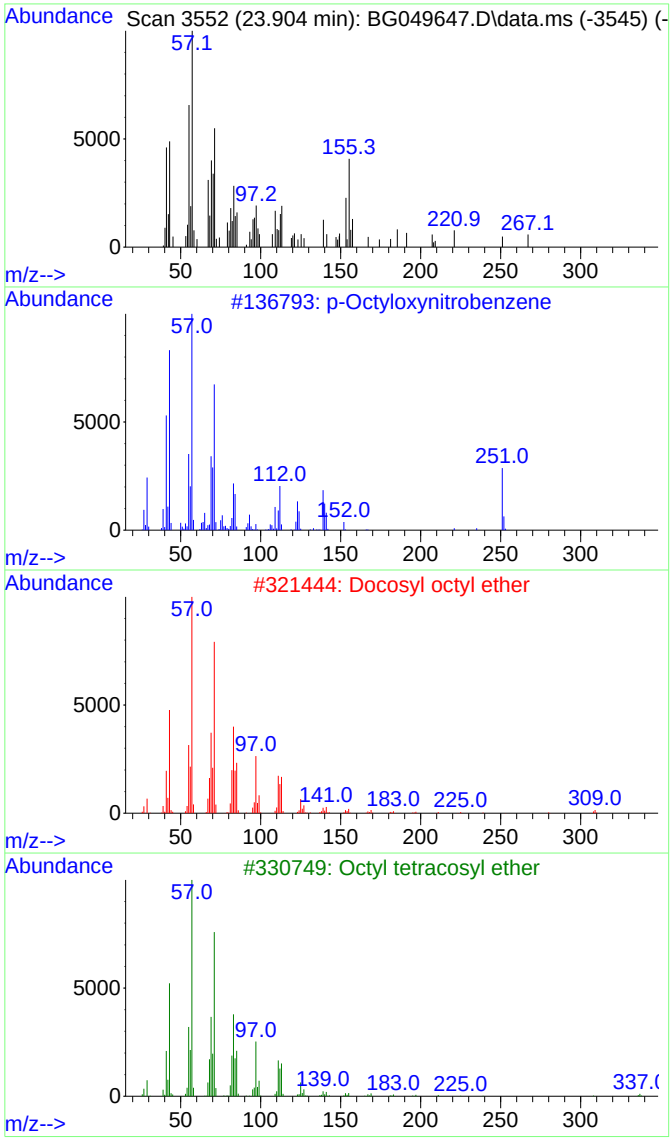
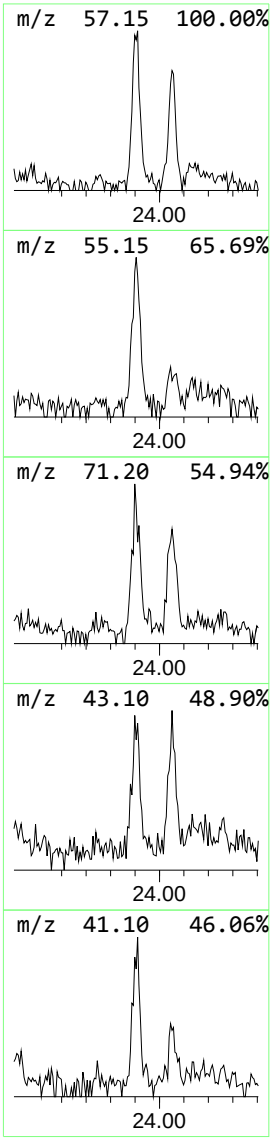
Peak Number 9 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.904	4.87 ng/ul	90856	Perylene-d12	25.003

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Octyloxynitrobenzene	251	C14H21NO3	049562-76-7	35
2	Docosyl octyl ether	438	C30H62O	1000406-38-9	30
3	Octyl tetracosyl ether	467	C32H66O	1000406-39-0	27
4	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	27
5	Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	22





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Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET3

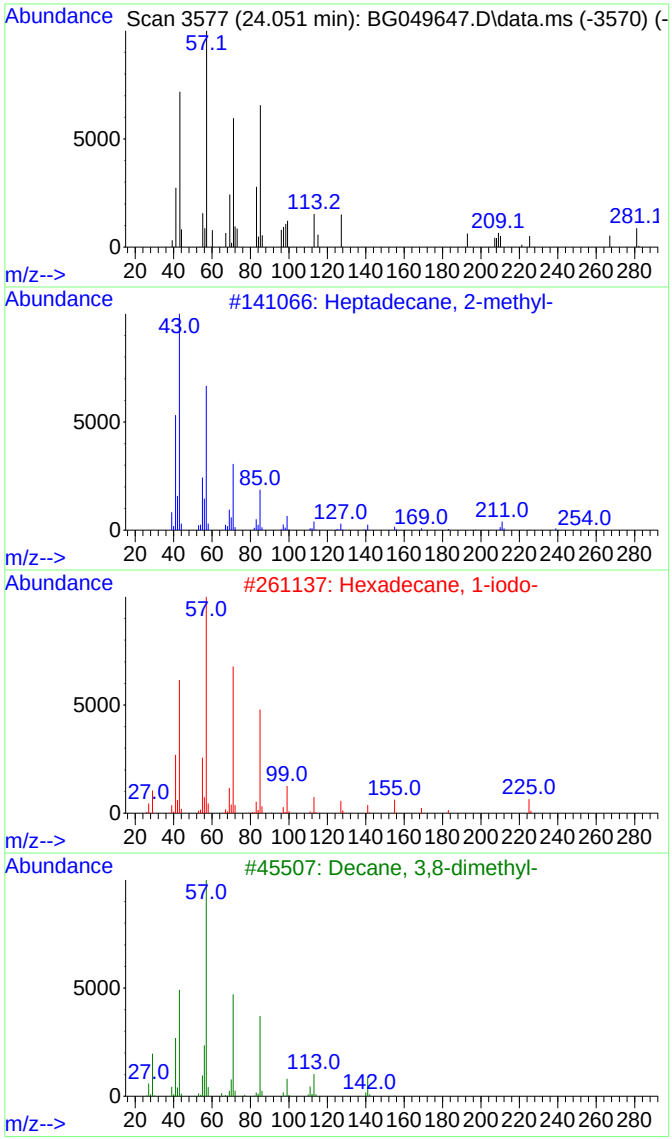
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

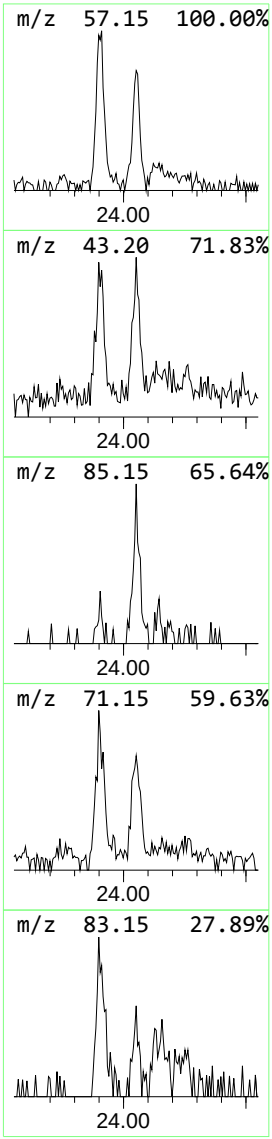
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.051	2.32 ng/ul	43386	Perylene-d12	25.003

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	59
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	59
3	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	53
4	Dotetracontane	591	C42H86	007098-20-6	53
5	Hentriacontane	437	C31H64	000630-04-6	50





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Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

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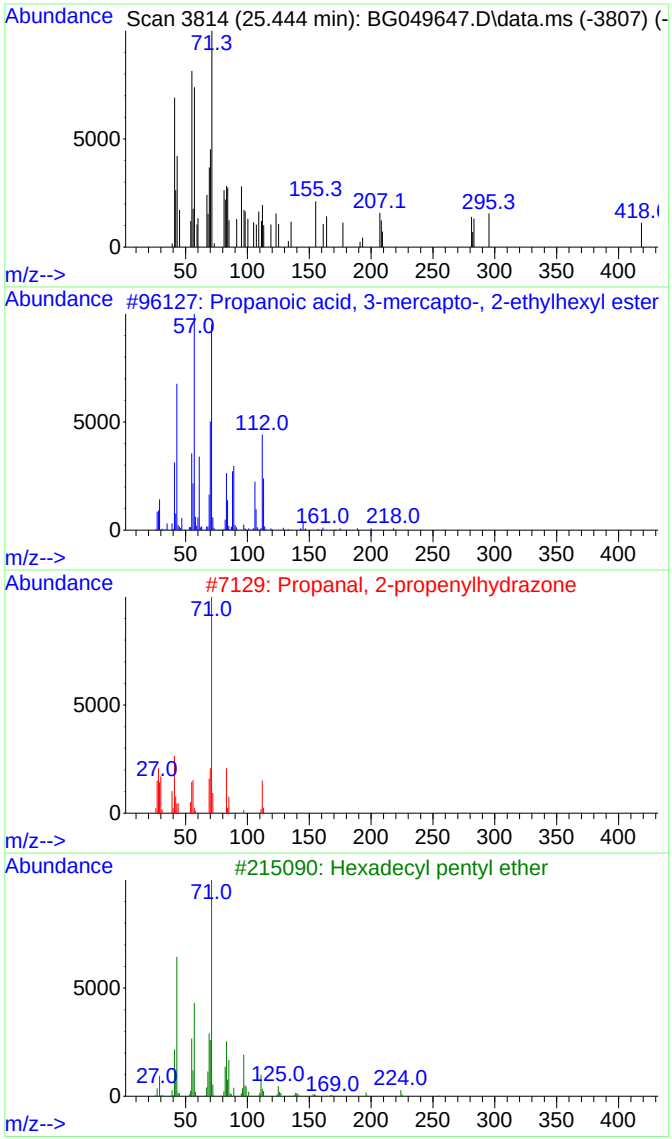
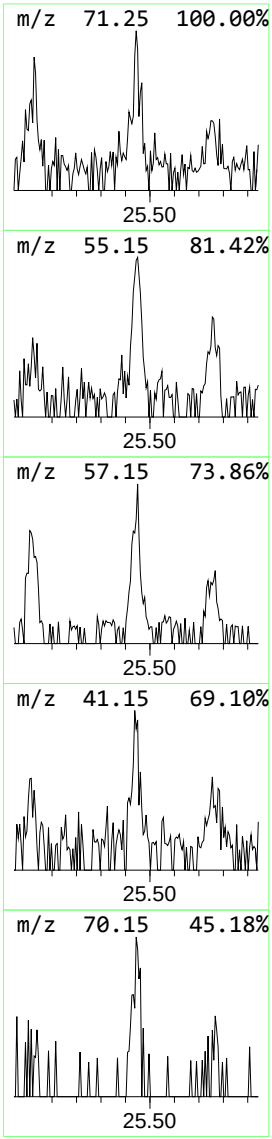
Peak Number 11 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.444	2.29 ng/ul	42717	Perylene-d12	25.003

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 3-mercapto-, 2-e...	218	C11H22O2S	050448-95-8	47
2	Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	38
3	Hexadecyl pentyl ether	312	C21H44O	1000406-41-8	35
4	Pentyl tetradecyl ether	284	C19H40O	1000406-41-7	35
5	Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	27





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049647.D
Acq On : 5 Aug 2021 13:12
Operator : CG/JU
Sample : M3162-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BGET3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.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...	3.074	17.2	ng/ul	120373	1	8.044	140006	20.0
Butane, 2-metho...	3.156	13.6	ng/ul	95509	1	8.044	140006	20.0
1-Octyn-3-ol, 4...	8.096	2.1	ng/ul	14513	1	8.044	140006	20.0
Ethanol, 2-(2-b...	10.628	4.7	ng/ul	46992	2	10.863	198659	20.0
1,3-Benzenediol...	18.759	11.3	ng/ul	187968	4	17.443	333462	20.0
Hexanedioic aci...	20.803	189.4	ng/ul	3238030	5	21.725	341931	20.0
Pentadecafluoro...	21.355	4.4	ng/ul	75578	5	21.725	341931	20.0
Tridecanol, 2-e...	22.530	2.3	ng/ul	39130	5	21.725	341931	20.0
unknown-01	23.904	4.9	ng/ul	90856	6	25.003	373234	20.0
(DEL) Alkane: S...	24.051	2.3	ng/ul	43386	6	25.003	373234	20.0
unknown-02	25.444	2.3	ng/ul	42717	6	25.003	373234	20.0