LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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: BG049282.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	7	13	23	rBV3	91491	232123	0.34%	0.128%
2	5.003	311	334	340	rBV3	74335	328122	0.48%	0.181%
3	5.162	342	361	369	rVV2	58035	216424	0.32%	0.120%
4	5.450	405	410	420	rBV	128657	209768	0.31%	0.116%
5	5.896	474	486	490	rBV	304540	601064	0.88%	0.332%
6	5.949	490	495	503	rVB	677815	1172411	1.72%	0.647%
7	6.102	516	521	537	rVB	225677	404380	0.59%	0.223%
8	7.154	694	700	705	rVV	114366	185163	0.27%	0.102%
9	7.271	705	720	732	rVV	2203679	4680637	6.88%	2.585%
10	7.606	771	777	782	rBV	88431	153359	0.23%	0.085%
11	7.718	790	796	803	rVV	229493	411025	0.60%	0.227%
12	7.847	803	818	841	rVB4	163578	907087	1.33%	0.501%
13	8.076	848	857	860	rBV	123176	228365	0.34%	0.126%
14	8.294	887	894	903	rBV	135334	235869	0.35%	0.130%
15	8.834	981	986	989	rVV	81880	146602	0.22%	0.081%
16	8.869	989	992	1003	rVB	77972	127757	0.19%	0.071%
17	9.239	1049	1055	1060	rVV3	83910	156998	0.23%	0.087%
18	9.304	1060	1066	1078	rVB	64802	147301	0.22%	0.081%
19	10.115	1197	1204	1213	rVB	242566	567854	0.83%	0.314%
20	10.362	1239	1246	1261	rVB	453795	788866	1.16%	0.436%
21	10.714	1280	1306	1317	rVB2	313329	1750352	2.57%	0.967%
22	10.896	1325	1337	1343	rBV2	206593	479247	0.70%	0.265%
23	11.267	1393	1400	1410	rBV2	147012	316744	0.47%	0.175%
24	11.690	1452	1472	1488	rBV	5462812	17770196	26.13%	9.812%
25	11.872	1488	1503	1520	rVB3	139436	778120	1.14%	0.430%
26	12.301	1565	1576	1584	rBV	81704	171122	0.25%	0.094%
27	12.512	1605	1612	1621	rVB	431480	758388	1.12%	0.419%
28	13.693	1800	1813	1823	rBV4	55203	165151	0.24%	0.091%
29	13.905	1842	1849	1853	rVV7	45027	136788	0.20%	0.076%
30	14.122	1879	1886	1890	rBV	140690	256407	0.38%	0.142%
31	14.169	1890	1894	1901	rVB	231943	400008	0.59%	0.221%
32	14.410	1923	1935	1941	rVB	176469	334090	0.49%	0.184%
33	14.651	1971	1976	1981	rVV	345138	465740	0.68%	0.257%
34	14.721	1981	1988	1993	rVV	265926	486180	0.71%	0.268%
35	14.921	2015	2022	2025	rBV2	86540	162319	0.24%	0.090%
36	14.956	2025	2028	2031	rVV	77964	127255	0.19%	0.070%

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Misc :
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Instrument : BNA_G
ClientSampleId : DBK43

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	14.997	2031	2035	2046	rVB3	209392	462081	0.68%	0.255%
38	15.121	2046	2056	2061	rBV	7028560	12097339	17.79%	6.680%
39	15.326	2082	2091	2100	rVB3	503896	948627	1.39%	0.524%
40	15.591	2112	2136	2140	rBV	16787108	68008730	100.00%	37.554%
41	15.750	2156	2163	2168	rVB	197109	353522	0.52%	0.195%
42	15.832	2169	2177	2182	rVV4	140362	296340	0.44%	0.164%
43	16.390	2268	2272	2280	rVB2	93120	136640	0.20%	0.075%
44	16.478	2281	2287	2291	rBV	102837	179212	0.26%	0.099%
45	16.537	2291	2297	2304	rBV	1246664	1807163	2.66%	0.998%
46	16.648	2309	2316	2324	rVV2	2123118	3791010	5.57%	2.093%
47	16.713	2324	2327	2331	rVV4	83180	140158	0.21%	0.077%
48	16.784	2331	2339	2343	rVV5	136101	378308	0.56%	0.209%
49	16.860	2347	2352	2361	rVV4	276001	582321	0.86%	0.322%
50	16.930	2361	2364	2373	rVV	105884	221771	0.33%	0.122%
51	17.195	2404	2409	2415	rVB	194757	306380	0.45%	0.169%
52	17.471	2450	2456	2462	rBV	325679	493996	0.73%	0.273%
53	17.571	2467	2473	2478	rVB	228735	366754	0.54%	0.203%
54	17.765	2500	2506	2513	rBV2	475545	757976	1.11%	0.419%
55	17.835	2513	2518	2525	rVB2	108152	192881	0.28%	0.107%
56	18.382	2601	2611	2623	rBV	2289178	4557848	6.70%	2.517%
57	18.693	2658	2664	2669	rBV	681399	1092266	1.61%	0.603%
58	18.775	2674	2678	2683	rBV7	78537	156013	0.23%	0.086%
59	19.198	2743	2750	2761	rVB	3184508	4530482	6.66%	2.502%
60	19.304	2763	2768	2773	rVB2	89002	134103	0.20%	0.074%
61	19.480	2793	2798	2801	rBV2	276078	452650	0.67%	0.250%
62	19.545	2805	2809	2817	rVV	395667	692453	1.02%	0.382%
63	19.639	2818	2825	2840	rVV	1769360	2835369	4.17%	1.566%
64	19.751	2840	2844	2848	rVB4	96635	126215	0.19%	0.070%
65	19.851	2857	2861	2867	rVV	267326	397422	0.58%	0.219%
66	19.915	2867	2872	2874	rVV2	367679	478597	0.70%	0.264%
67	19.950	2874	2878	2882	rVV	930360	1296326	1.91%	0.716%
68	19.992	2882	2885	2889	rVV	268198	359155	0.53%	0.198%
69	20.039	2889	2893	2898	rVV2	271951	382702	0.56%	0.211%
70	20.232	2920	2926	2932	rVB	407184	592588	0.87%	0.327%
71	20.338	2938	2944	2949	rBV3	136155	226642	0.33%	0.125%
72	20.597	2982	2988	2991	rBV2	71538	183788	0.27%	0.101%
73	20.708	2999	3007	3010	rBV3	148049	311043	0.46%	0.172%
74	21.131	3074	3079	3083	rBV2	95509	154081	0.23%	0.085%
75	21.178	3083	3087	3093	rVB	469321	677074	1.00%	0.374%
76	21.355	3107	3117	3136	rVB	4678983	8816537	12.96%	4.868%

LSC Area Percent Report

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Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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

77	21.519	3140	3145	3153	rVB3	132569	221967	0.33%	0.123%
78	21.719	3171	3179	3183	rBV	2172702	3587939	5.28%	1.981%
79	21.754	3183	3185	3193	rVB	326250	465429	0.68%	0.257%
80	22.465	3300	3306	3317	rVB6	229645	580543	0.85%	0.321%
81	23.247	3433	3439	3446	rVB	198941	400757	0.59%	0.221%
82	23.352	3452	3457	3463	rBV4	97221	206757	0.30%	0.114%
83	23.417	3463	3468	3472	rVV3	74792	146515	0.22%	0.081%
84	23.758	3518	3526	3540	rBV3	240822	600730	0.88%	0.332%
85	24.751	3688	3695	3701	rBV8	89334	280893	0.41%	0.155%
86	24.815	3702	3706	3716	rVB3	132959	330849	0.49%	0.183%
87	25.045	3736	3745	3756	rVB2	189074	567166	0.83%	0.313%
88	25.385	3795	3803	3815	rVB2	167323	485712	0.71%	0.268%
89	25.532	3820	3828	3835	rBV3	80461	264831	0.39%	0.146%
90	25.838	3871	3880	3881	rBV2	66144	167511	0.25%	0.092%
91	26.149	3925	3933	3939	rBV	150673	460584	0.68%	0.254%
92	26.231	3940	3947	3961	rVB2	279347	914989	1.35%	0.505%
93	26.754	4029	4036	4037	rBV6	94250	147429	0.22%	0.081%
94	26.860	4043	4054	4086	rVB3	919080	3834376	5.64%	2.117%
95	27.348	4118	4137	4158	rBV3	1651630	7529081	11.07%	4.157%
96	28.958	4405	4411	4428	rVB3	63378	287201	0.42%	0.159%
97	29.198	4438	4452	4466	rBV4	79139	351895	0.52%	0.194%
98	29.375	4469	4482	4499	rBV4	70873	390239	0.57%	0.215%
99	29.874	4551	4567	4588	rBV4	129427	754576	1.11%	0.417%
100	30.538	4664	4680	4712	rVB8	260469	1685799	2.48%	0.931%

Sum of corrected areas: 181097613

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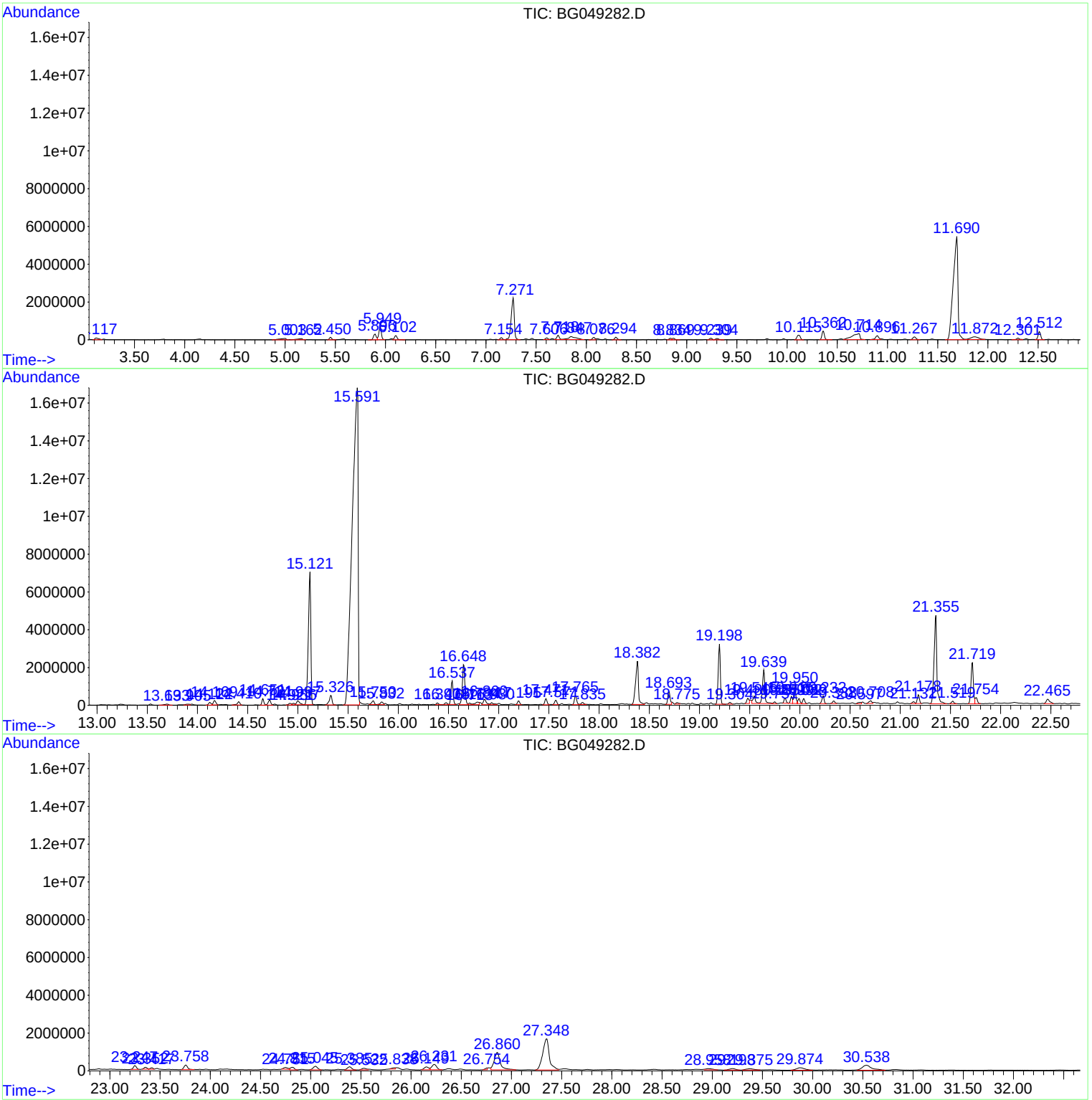
Instrument :
BNA_G
ClientSampleId :
DBK43

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

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



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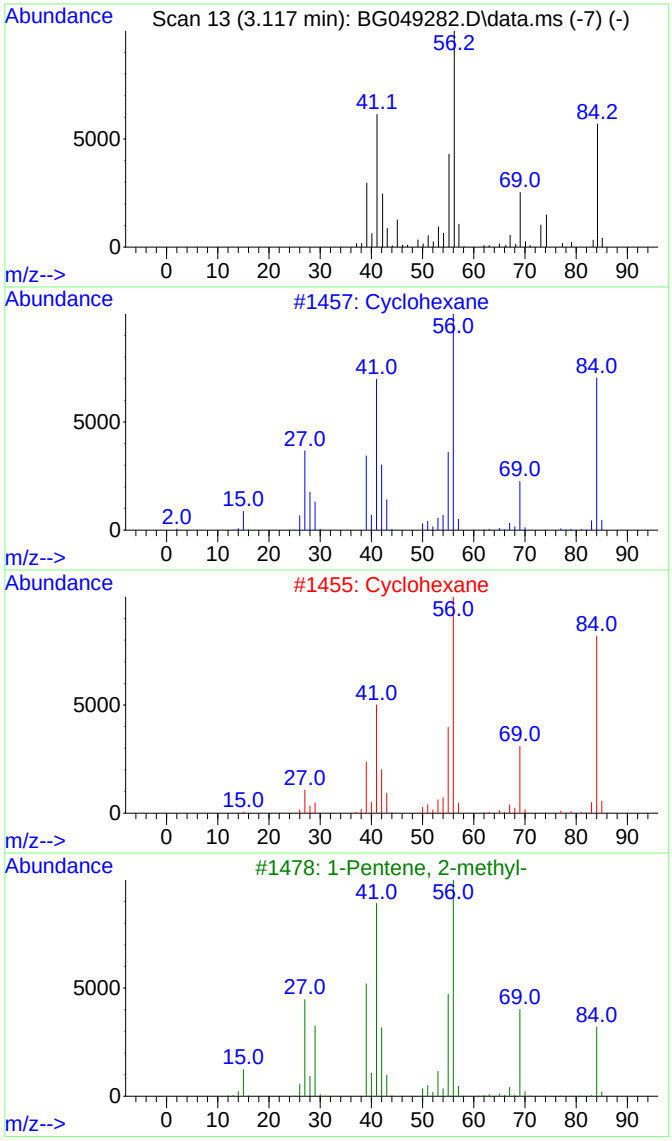
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

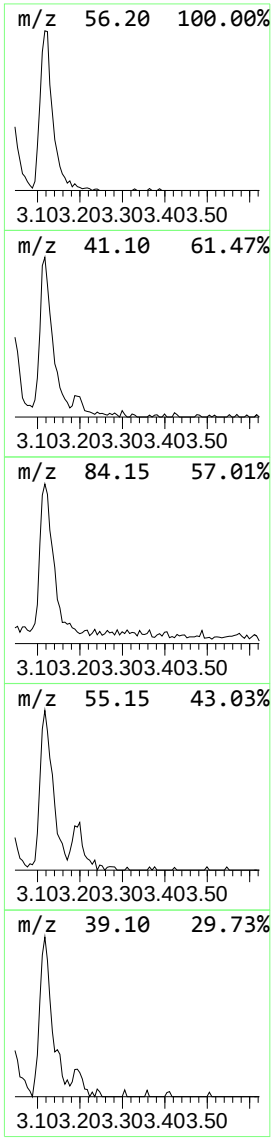
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.12 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.120	20.33 ng/ul	232123	1,4-Dichlorobenzene-d4	8.076

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	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	83
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80
5	Cyclohexane	84	C6H12	000110-82-7	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
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Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

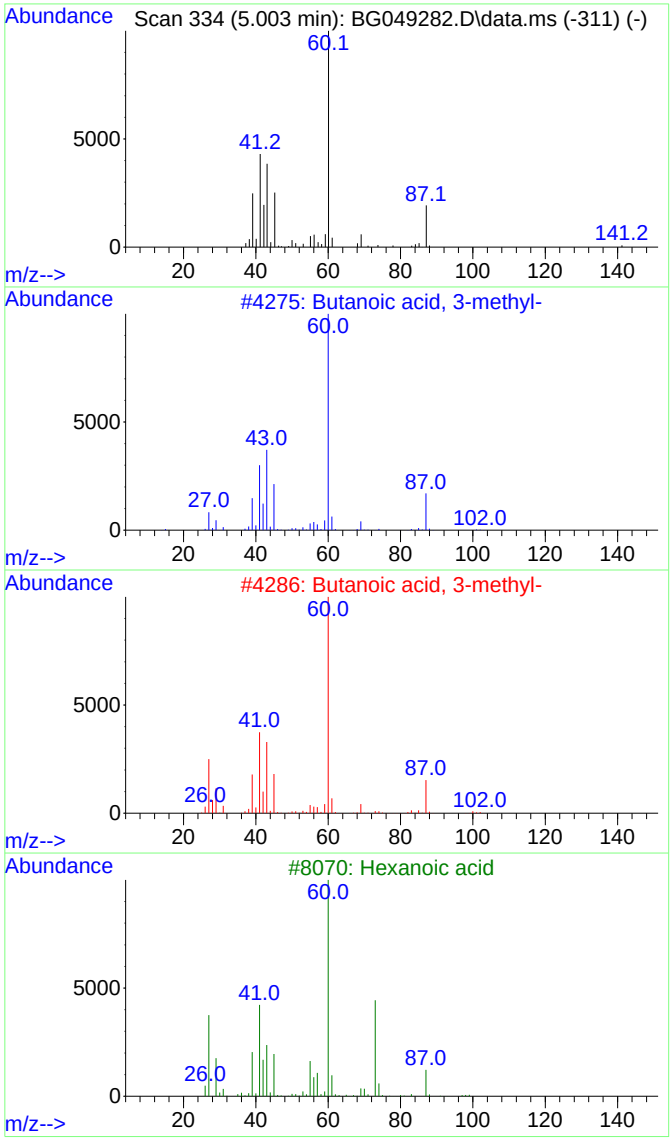
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

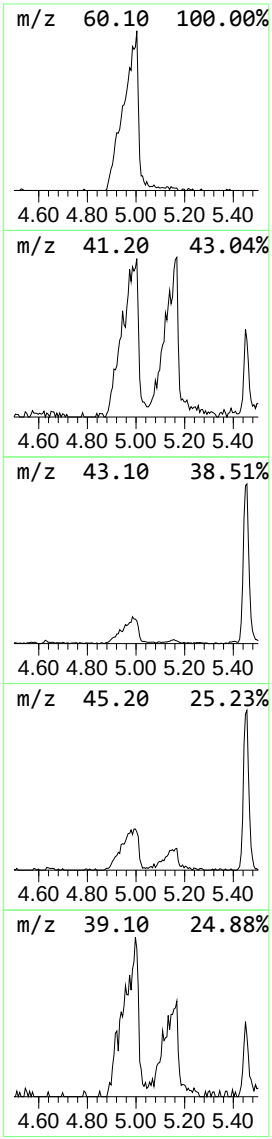
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butanoic acid, 3-methyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.000	28.74 ng/ul	328122	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	72
2	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
3	Hexanoic acid	116	C6H12O2	000142-62-1	39
4	Formic acid hydrazide	60	CH4N2O	000624-84-0	33
5	Pentanoic acid	102	C5H10O2	000109-52-4	28





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Instrument : BNA_G
ClientSampleId : DBK43

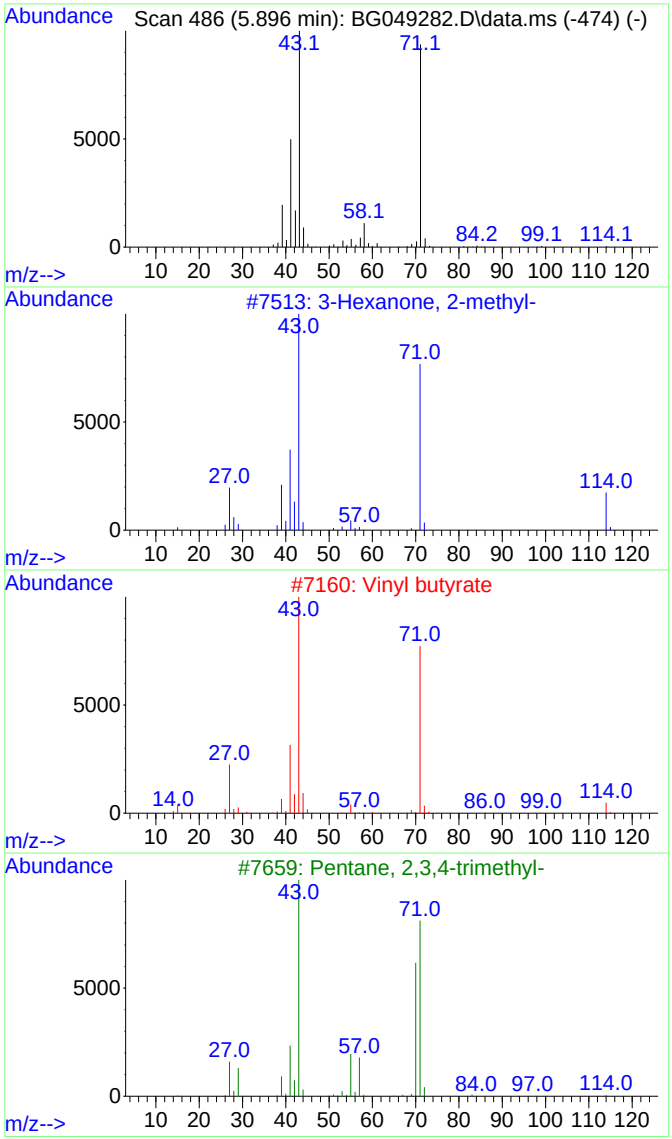
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

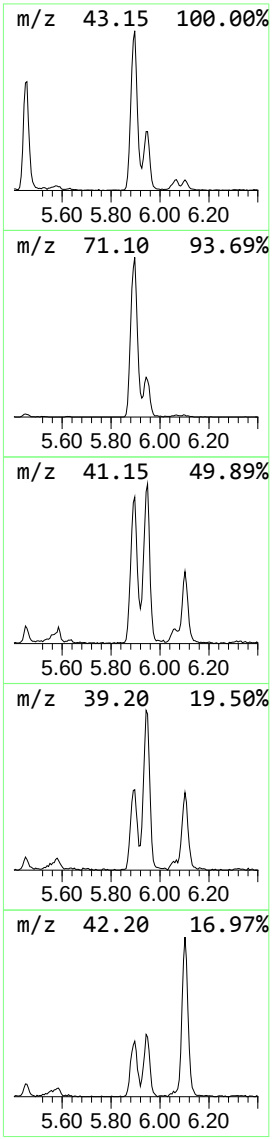
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 3-Hexanone, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.900	52.64 ng/ul	601064	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone, 2-methyl-	114	C7H14O	007379-12-6	64
2	Vinyl butyrate	114	C6H10O2	000123-20-6	64
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
4	Tetrahydro[2,2']bifuranyl-5-one	156	C8H12O3	019680-00-3	50
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	46





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Instrument : BNA_G
ClientSampleId : DBK43

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

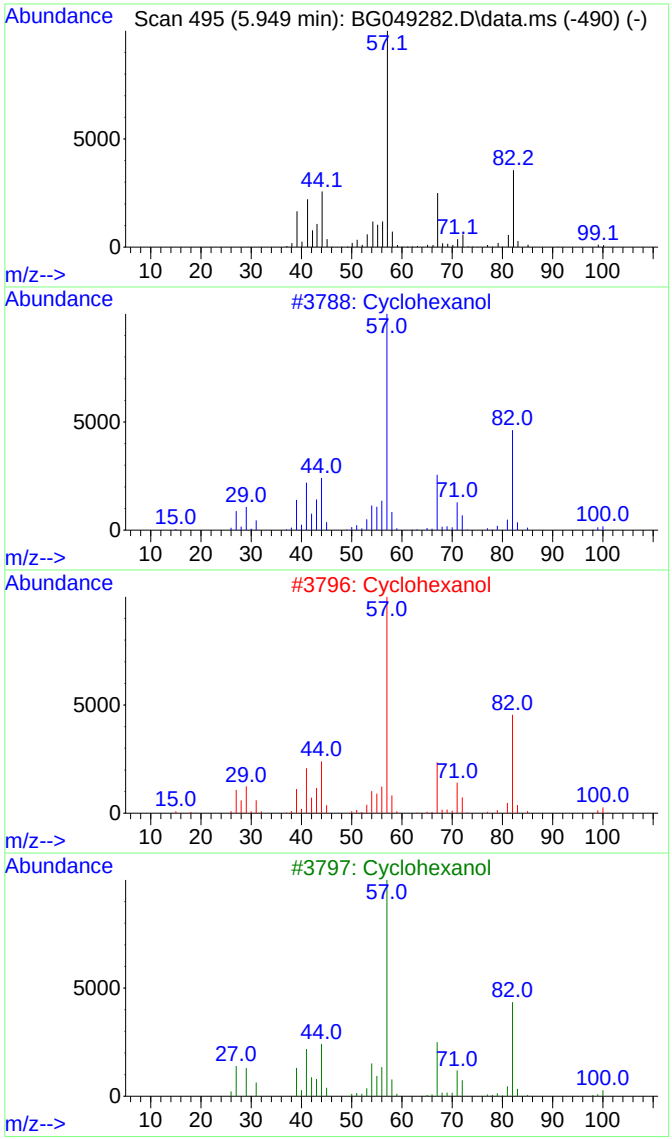
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

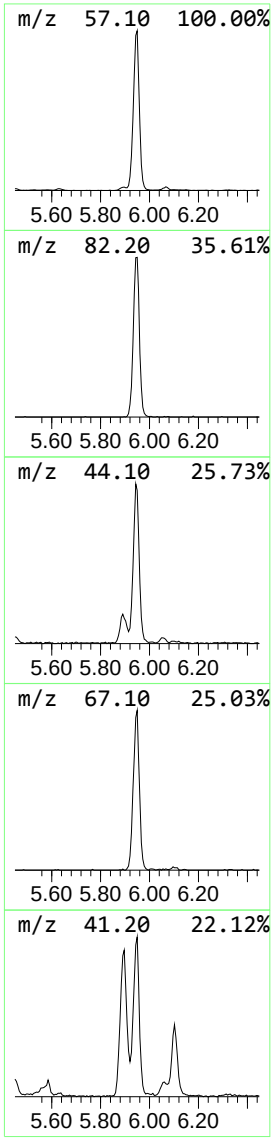
****

Peak Number 6 Cyclohexanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.950	102.68 ng/ul	1172410	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol	100	C6H12O	000108-93-0	91
2	Cyclohexanol	100	C6H12O	000108-93-0	83
3	Cyclohexanol	100	C6H12O	000108-93-0	83
4	Cyclohexanol	100	C6H12O	000108-93-0	64
5	2-Hexen-1-ol, (Z)-	100	C6H12O	000928-94-9	59





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Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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

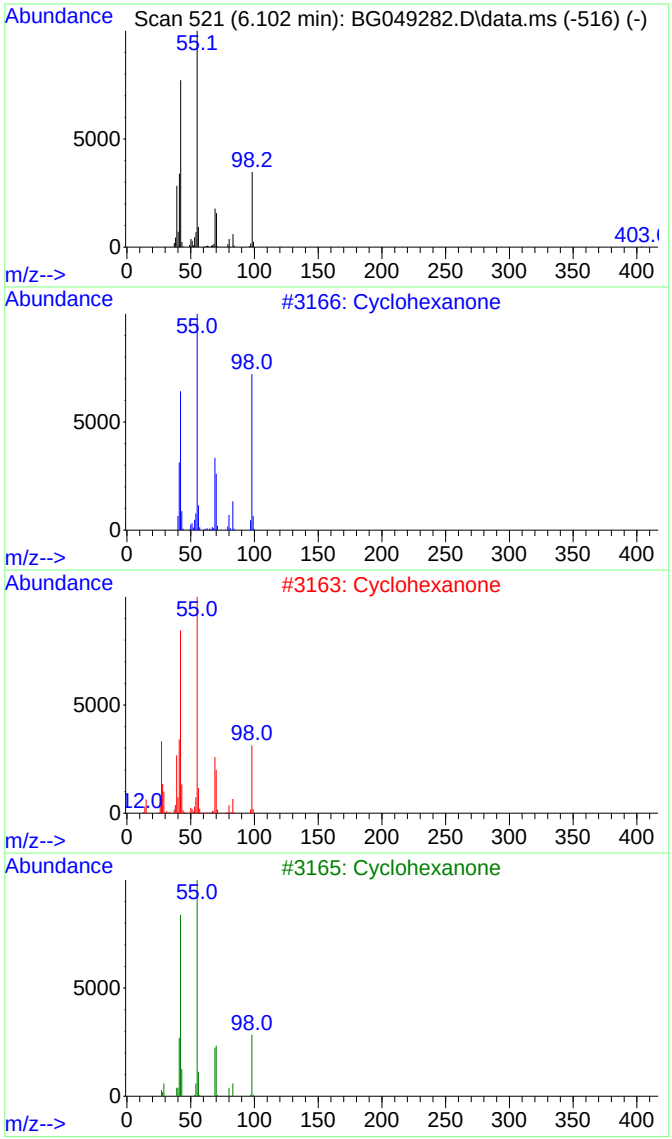
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

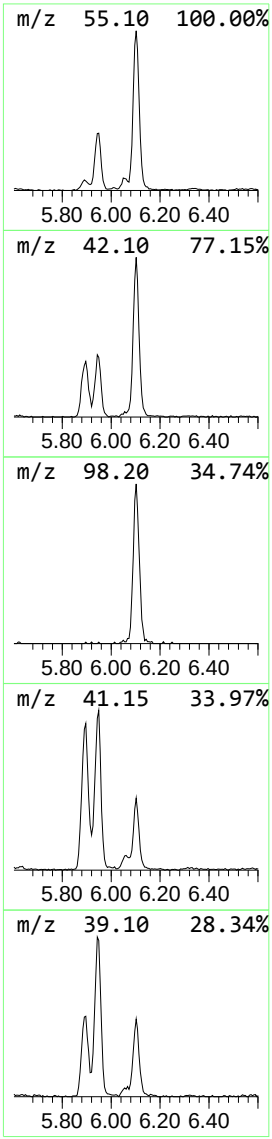
****

Peak Number 7 Cyclohexanone Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.100	35.42 ng/ul	404380	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone	98	C6H10O	000108-94-1	90
2	Cyclohexanone	98	C6H10O	000108-94-1	72
3	Cyclohexanone	98	C6H10O	000108-94-1	64
4	Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	64
5	Cyclohexanone	98	C6H10O	000108-94-1	62





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Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

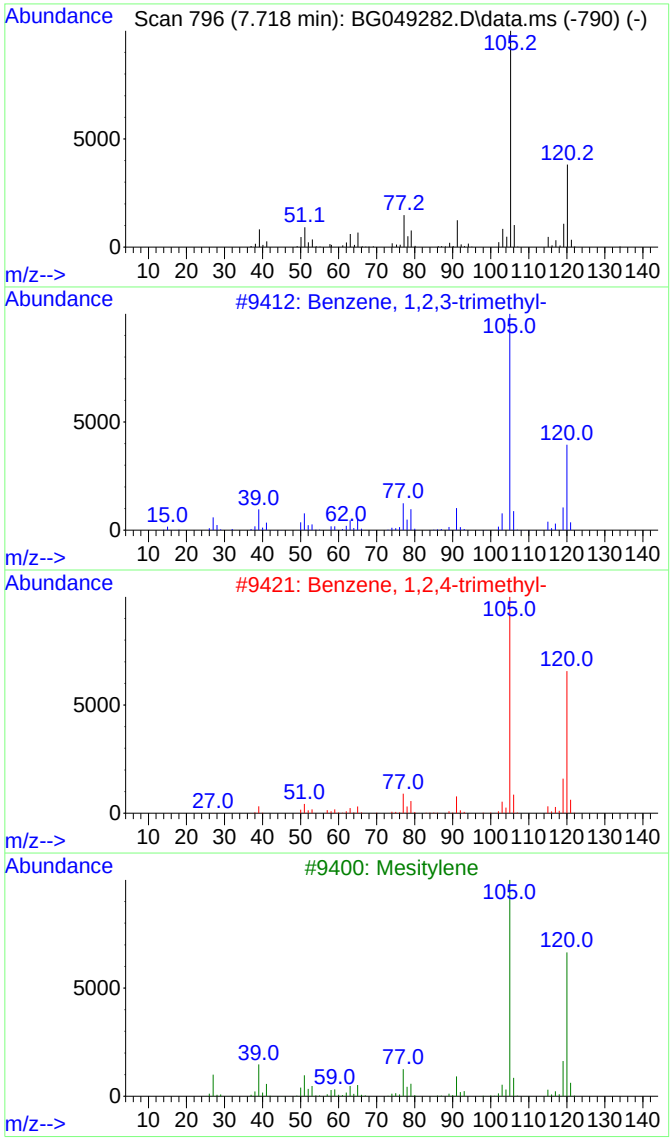
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

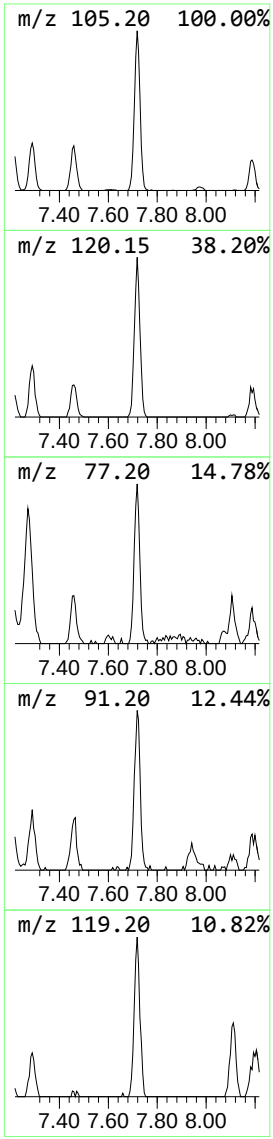
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.720	36.00 ng/ul	411025	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





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Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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

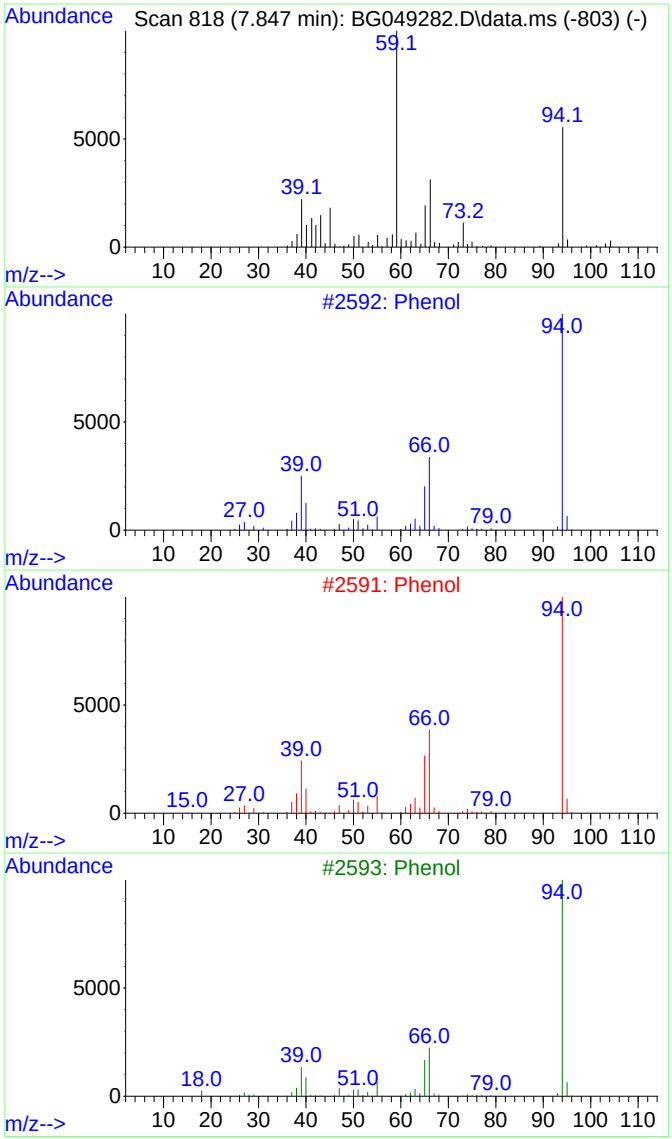
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

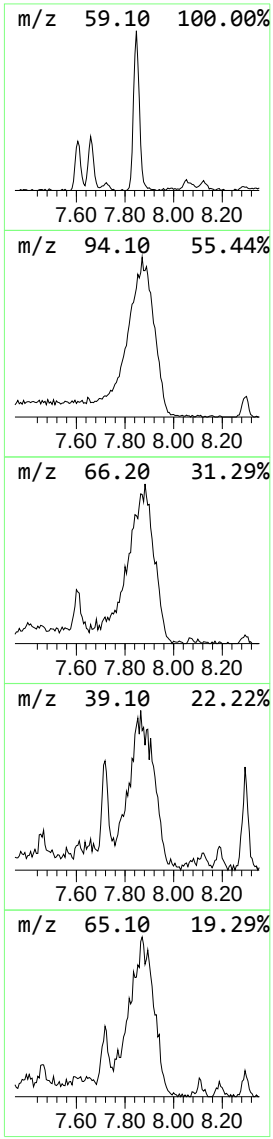
****

Peak Number 10 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.850	79.44 ng/ul	907087	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol	94	C6H6O	000108-95-2	60
2	Phenol	94	C6H6O	000108-95-2	60
3	Phenol	94	C6H6O	000108-95-2	46
4	Phosphonic acid, (p-hydroxyphenyl)-	174	C6H7O4P	033795-18-5	43
5	Carbamic acid, phenyl ester	137	C7H7NO2	000622-46-8	38





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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

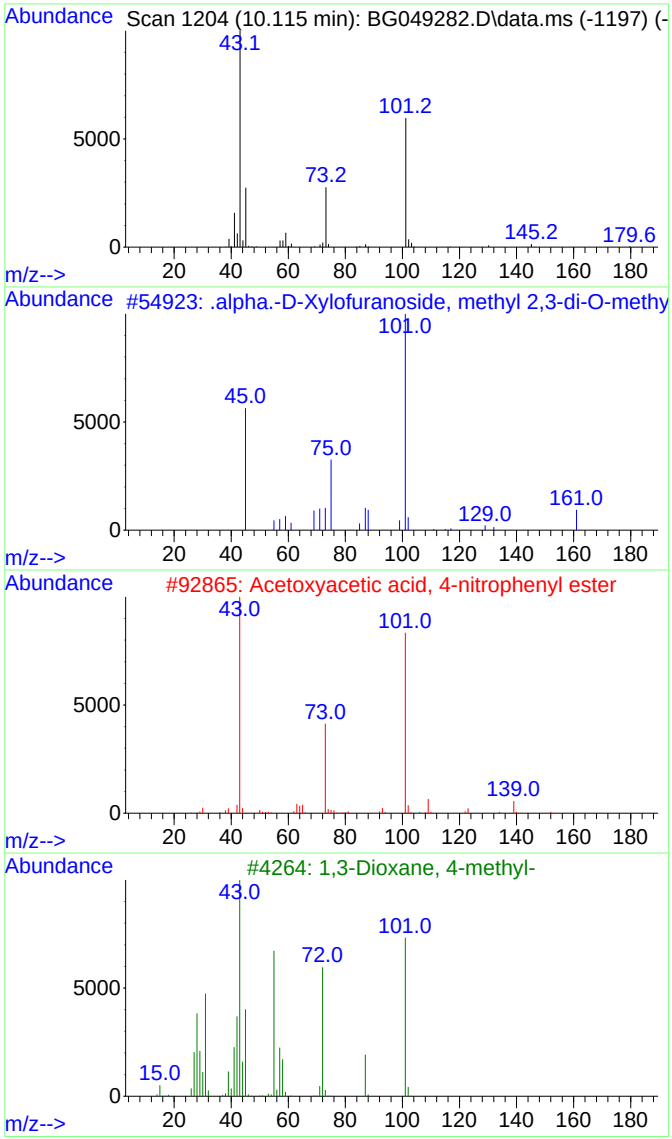
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

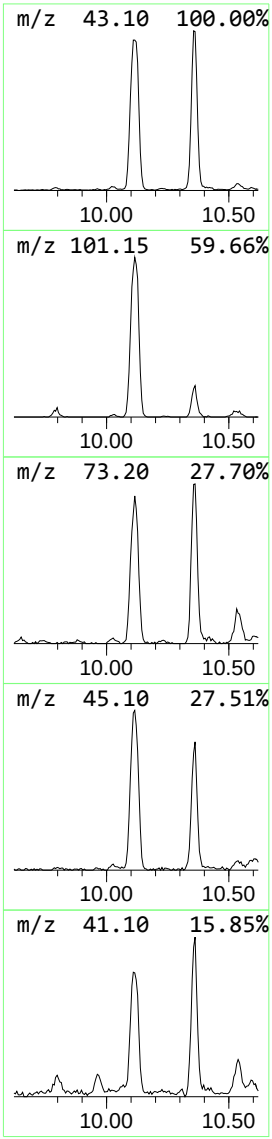
****

Peak Number 13 unknown-02 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.120	23.70 ng/ul	567854	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-D-Xylofuranoside, methyl...	192	C8H16O5	015821-56-4	38
2	Acetoxyacetic acid, 4-nitropheny...	239	C10H9NO6	1000307-54-5	36
3	1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	36
4	Hexahydro-5H-imidazo[5,1-b:4,3-b...	188	C7H12N2S2	019505-80-7	33
5	Acetoxyacetic acid, 3-methylbut-...	188	C9H16O4	1000355-69-8	33





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

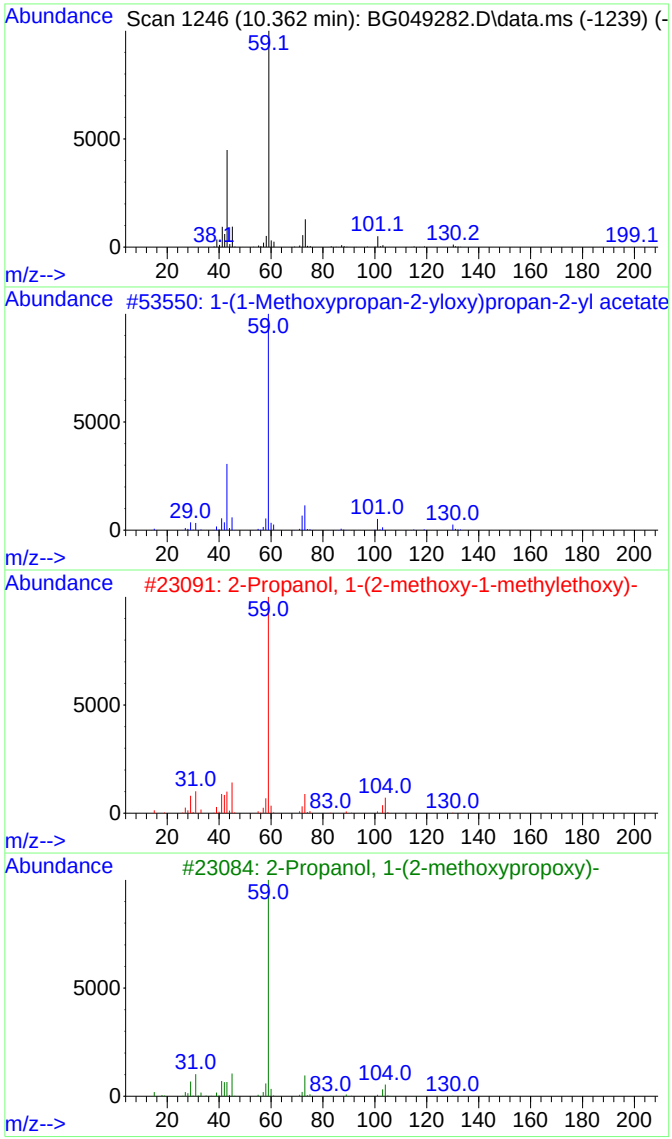
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

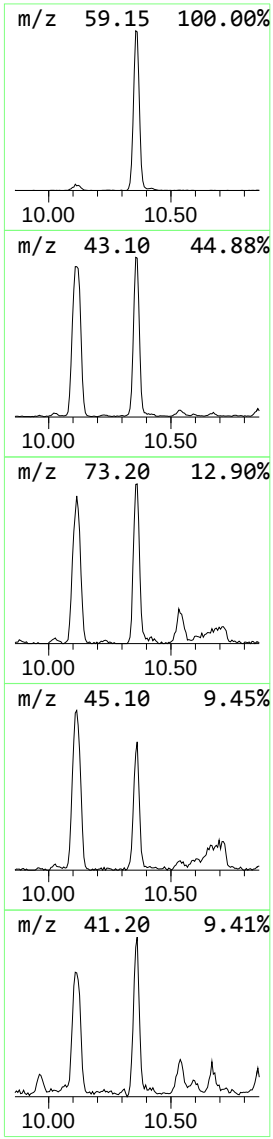
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	32.92 ng/ul	788866	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	83
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
3	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	72
4	Amylene hydrate	88	C5H12O	000075-85-4	64
5	1-(1-Methoxypropan-2-yloxy)propa...	244	C9H15F3O4	1000367-08-7	64





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Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

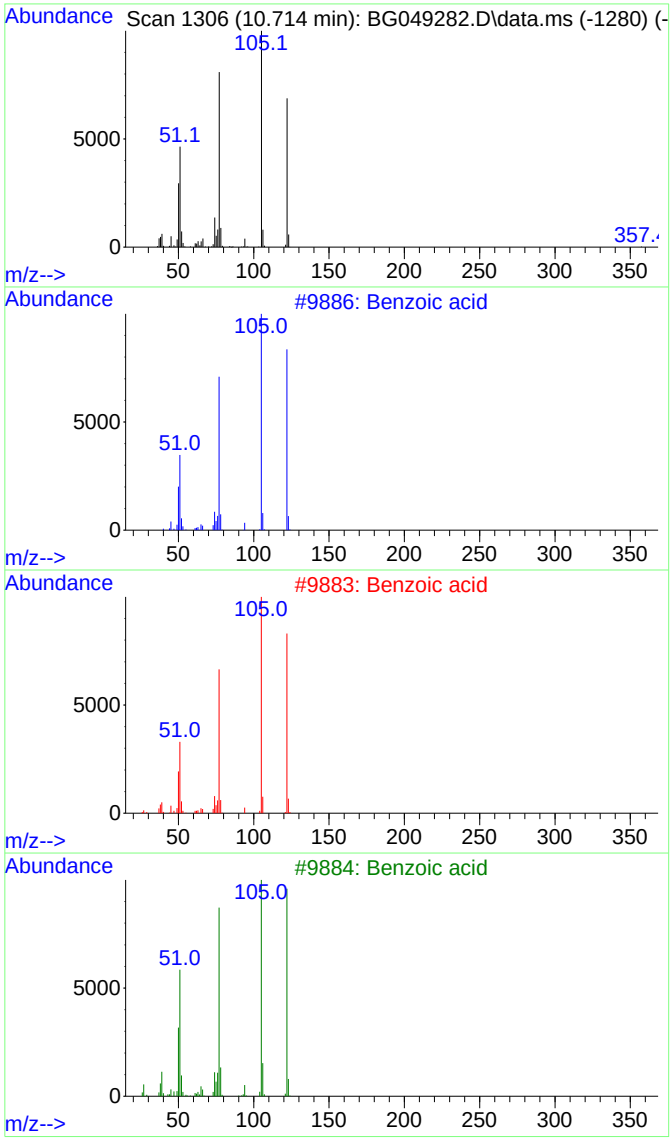
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

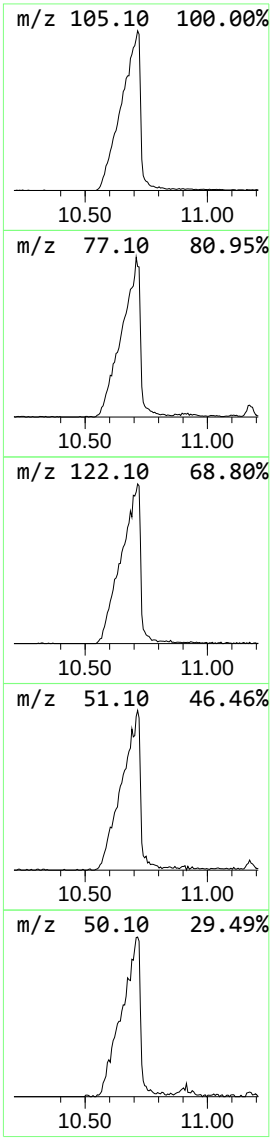
****

Peak Number 15 Benzoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.710	73.05 ng/ul	1750350	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	96
2	Benzoic acid	122	C7H6O2	000065-85-0	94
3	Benzoic acid	122	C7H6O2	000065-85-0	90
4	Benzoic acid, silver(1+) salt	228	C7H5AgO2	000532-31-0	90
5	Cyclobutane-1,1-dicarboxamide, N...	382	C20H18N2O6	1000253-25-3	90





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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

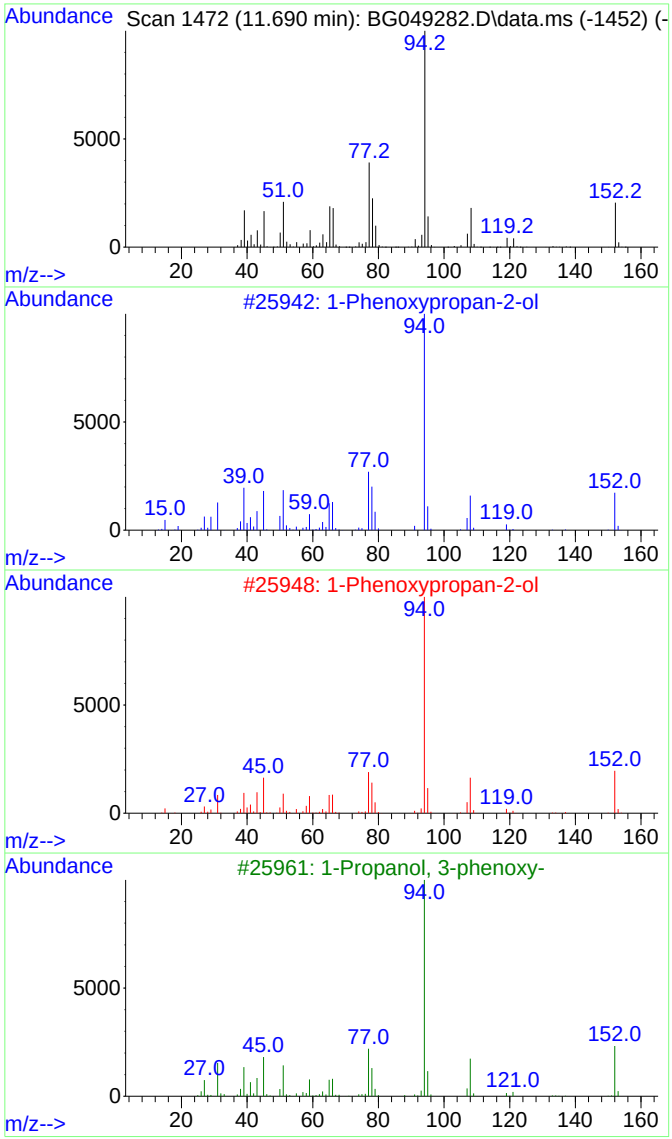
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

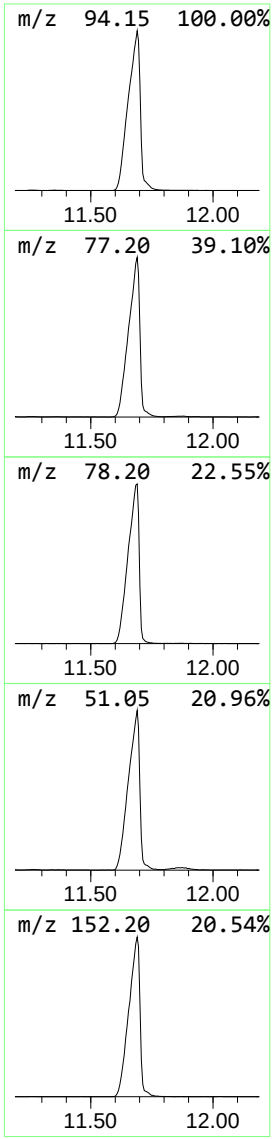
****

Peak Number 17 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.690	741.59 ng/ul	17770200	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	92
3	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	64
4	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	59
5	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	59





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Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

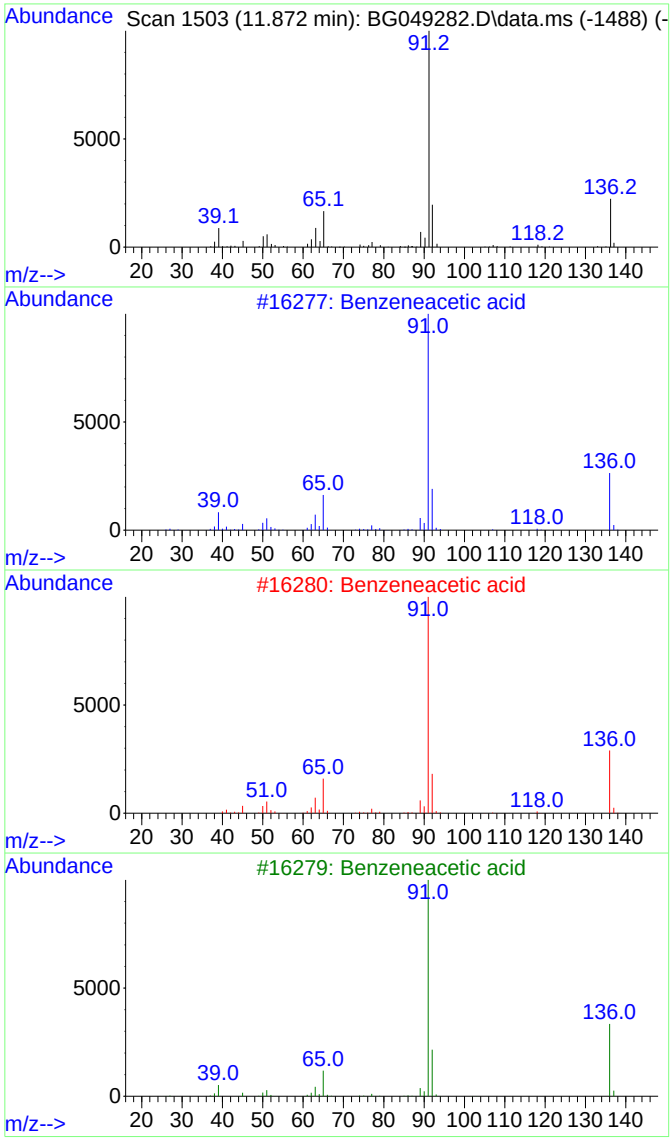
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

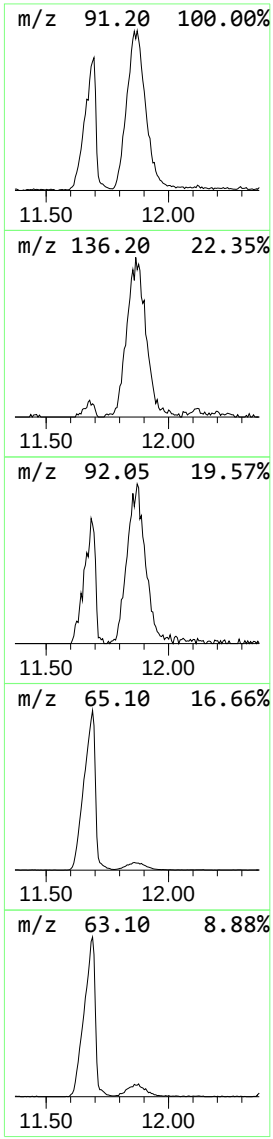
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzeneacetic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.870	32.47 ng/ul	778120	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
2	Benzeneacetic acid	136	C8H8O2	000103-82-2	91
3	Benzeneacetic acid	136	C8H8O2	000103-82-2	90
4	Propanedioic acid, phenyl-	180	C9H8O4	002613-89-0	83
5	Acetic acid, phenyl-, isopentyl ...	206	C13H18O2	000102-19-2	64





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ALS Vial : 53 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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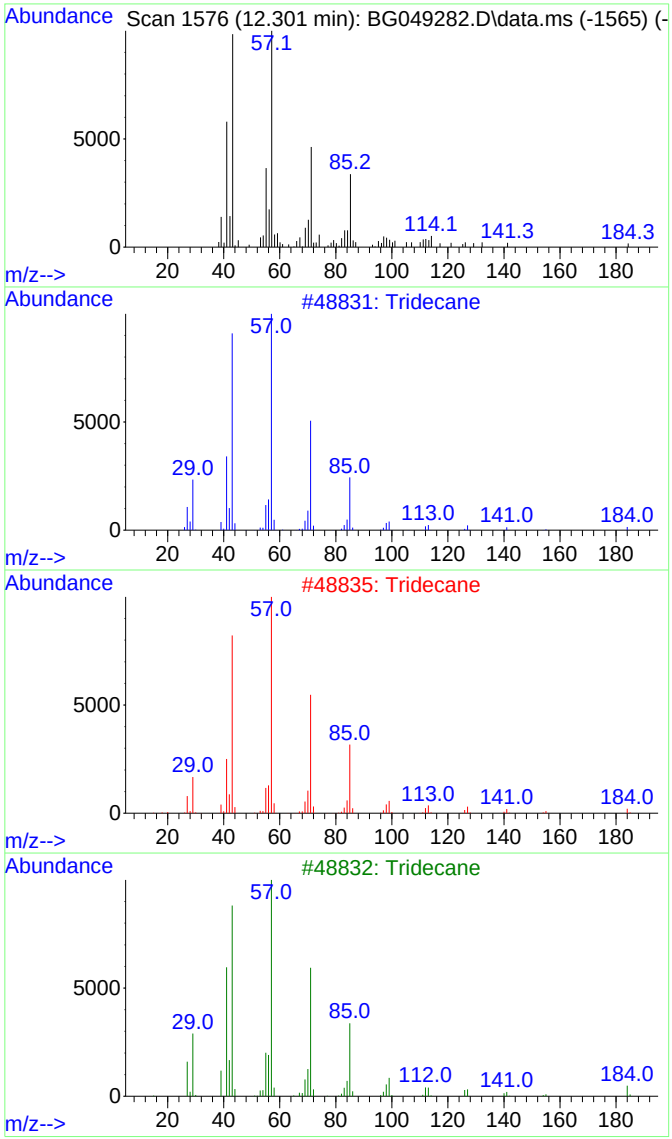
Quant Title : SVOA CALIBRATION

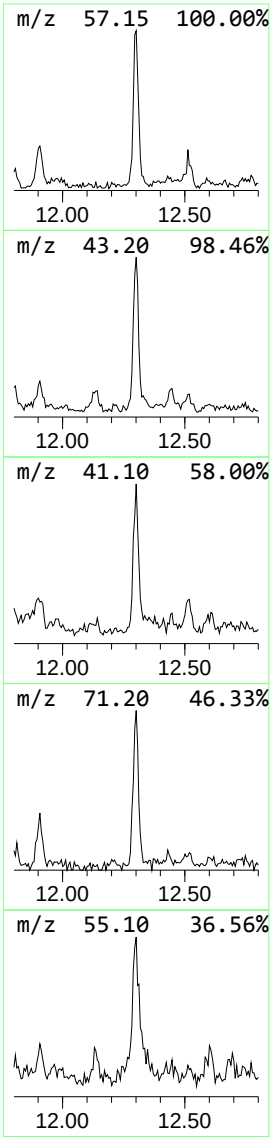
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.300	7.14 ng/ul	171122	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	87
2	Tridecane	184	C13H28	000629-50-5	64
3	Tridecane	184	C13H28	000629-50-5	64
4	Hexadecane	226	C16H34	000544-76-3	64
5	Heptadecane	240	C17H36	000629-78-7	59





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Sample : M2914-14
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Instrument : BNA_G
ClientSampleId : DBK43

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Quant Title : SVOA CALIBRATION

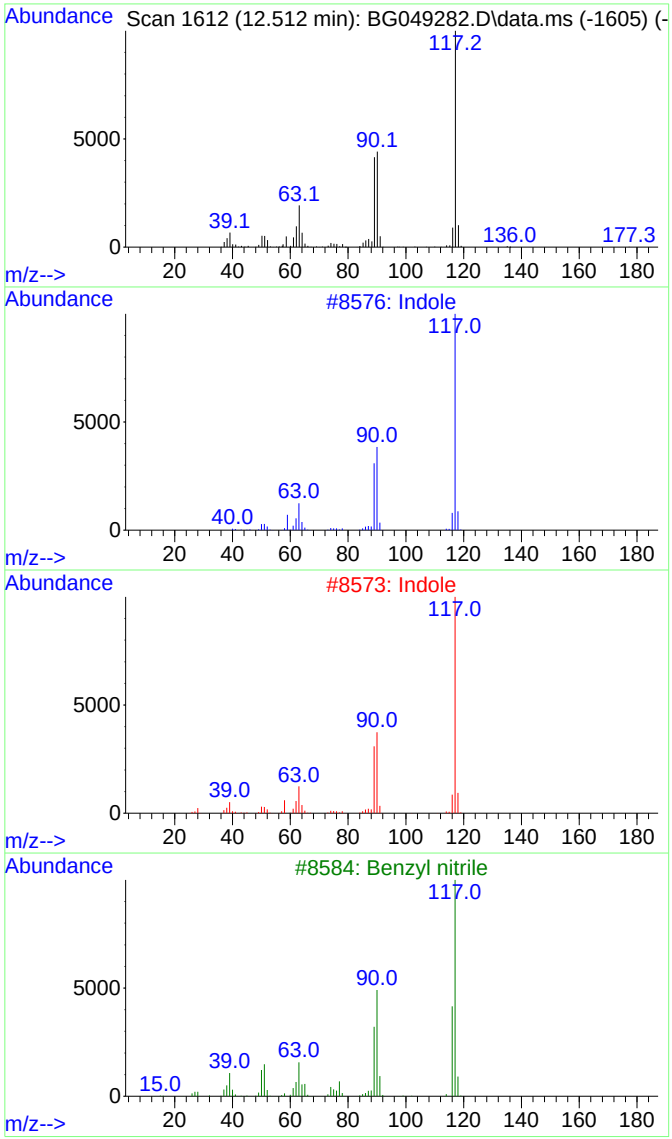
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

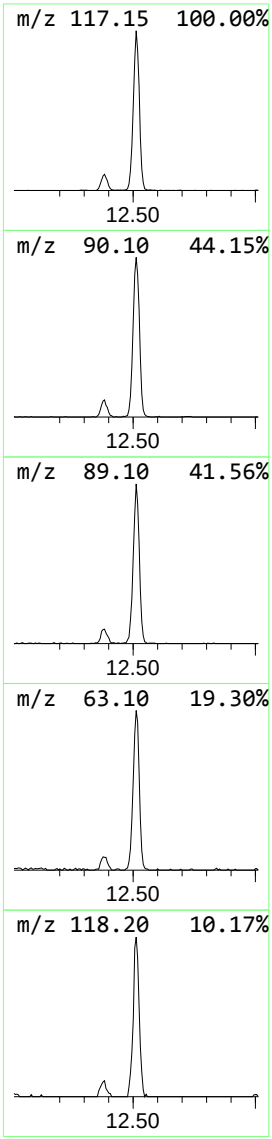
****

Peak Number 20 Indole Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.510	31.65 ng/ul	758388	Naphthalene-d8	10.896

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole	117	C8H7N	000120-72-9	95
2	Indole	117	C8H7N	000120-72-9	94
3	Benzyl nitrile	117	C8H7N	000140-29-4	91
4	Indolizine	117	C8H7N	000274-40-8	91
5	Indole	117	C8H7N	000120-72-9	90





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Acq On : 11 Jul 2021 2:38
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Sample : M2914-14
Misc :
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Instrument : BNA_G
ClientSampleId : DBK43

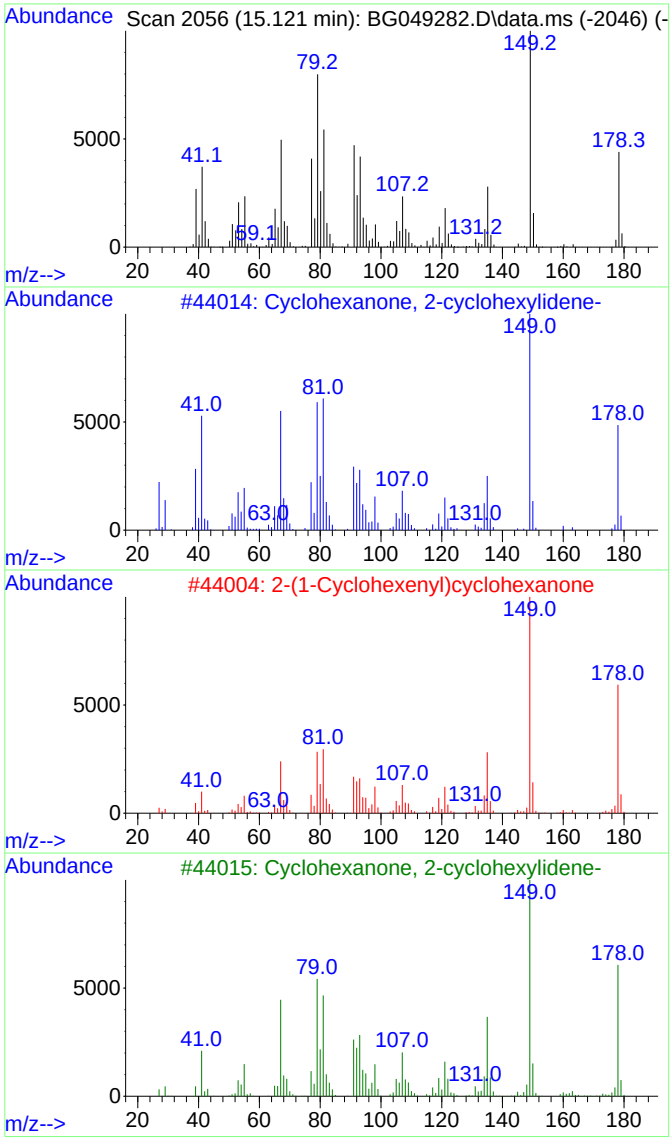
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

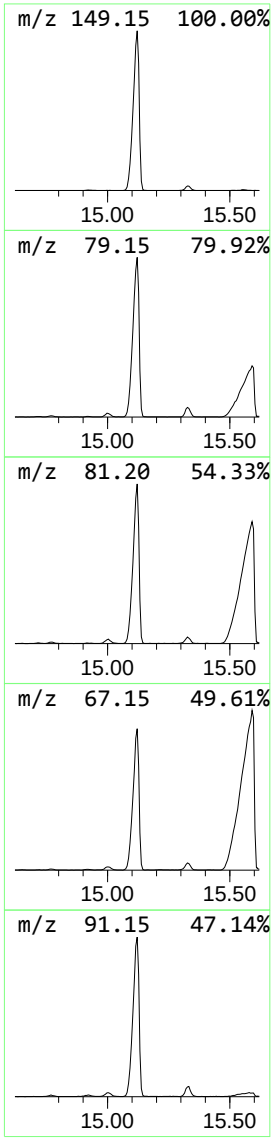
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 Cyclohexanone, 2-cyclohexyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.120	497.65 ng/ul	12097300	Acenaphthene-d10	14.721

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	93
2	2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	92
3	Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	60
4	Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	55
5	3-Nonen-1-yne, (Z)-	122	C9H14	037981-61-6	50





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Sample : M2914-14
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ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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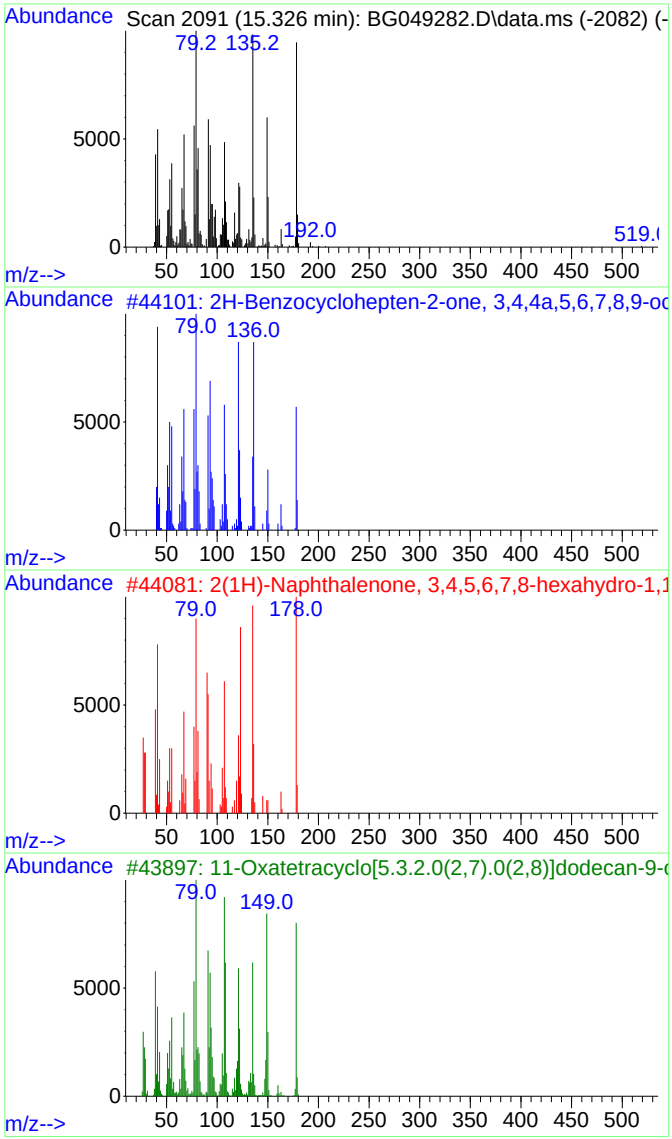
Quant Title : SVOA CALIBRATION

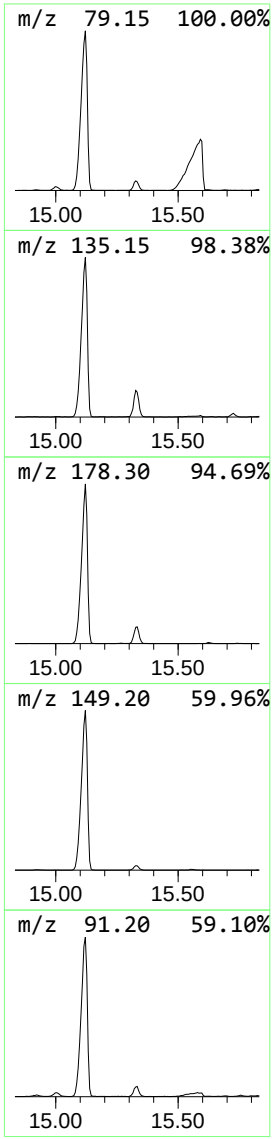
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 2H-Benzocyclohepten-2-one, ... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.330	39.02 ng/ul	948627	Acenaphthene-d10	14.721

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Benzocyclohepten-2-one, 3,4,4...	178	C12H18O	055103-71-4	70
2	2(1H)-Naphthalenone, 3,4,5,6,7,8...	178	C12H18O	001609-25-2	64
3	11-Oxatetracyclo[5.3.2.0(2,7).0(...	178	C11H14O2	1000186-25-4	62
4	p-Cymen-7-ol	150	C10H14O	000536-60-7	55
5	Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	53





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Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

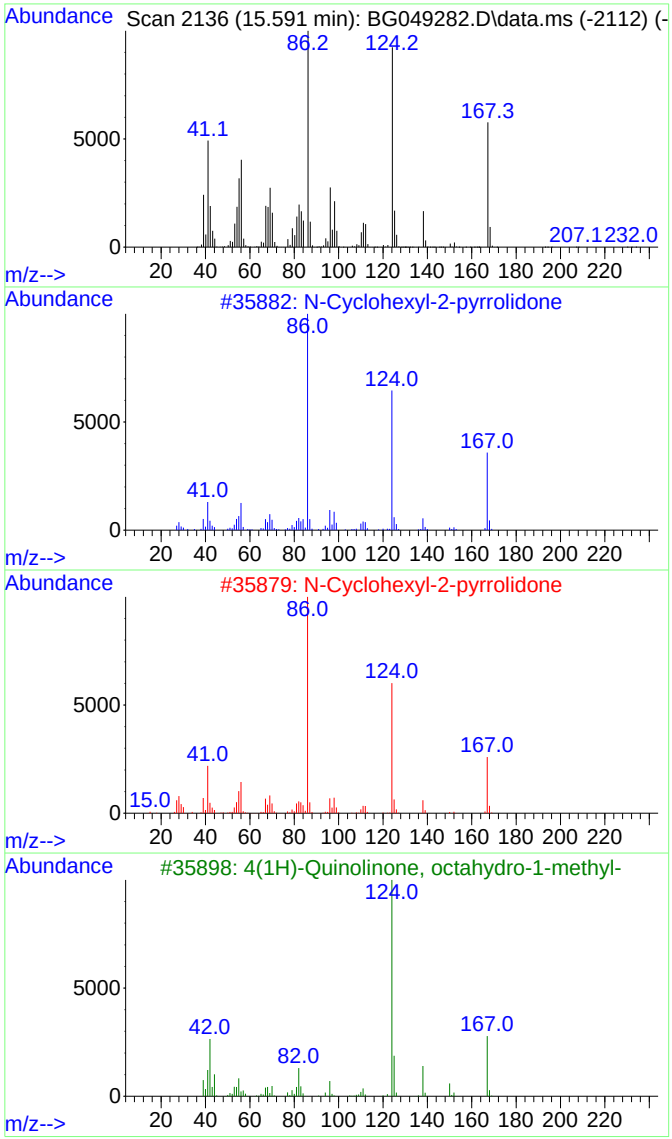
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

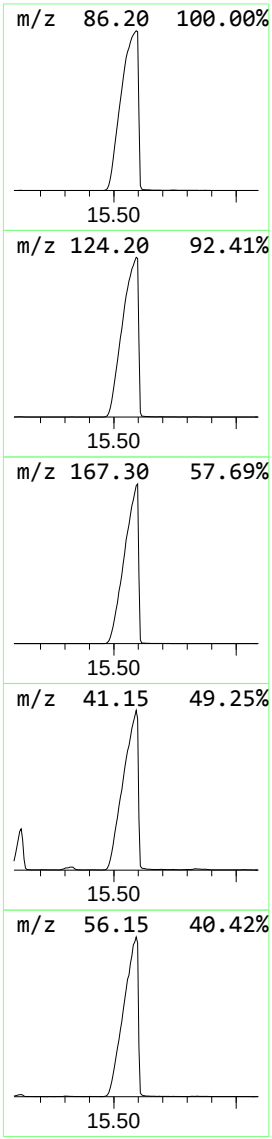
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 N-Cyclohexyl-2-pyrrolidone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.590	2797.68 ng/ul	68008700	Acenaphthene-d10	14.721

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Cyclohexyl-2-pyrrolidone	167	C10H17NO	006837-24-7	53
2	N-Cyclohexyl-2-pyrrolidone	167	C10H17NO	006837-24-7	38
3	4(1H)-Quinolinone, octahydro-1-m...	167	C10H17NO	1000320-16-7	35
4	Piperidine, 1-(2-methyl-1-propen...	139	C9H17N	000673-33-6	16
5	4,6-Dimethyl-2-pyrimidone	124	C6H8N2O	000108-79-2	14





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

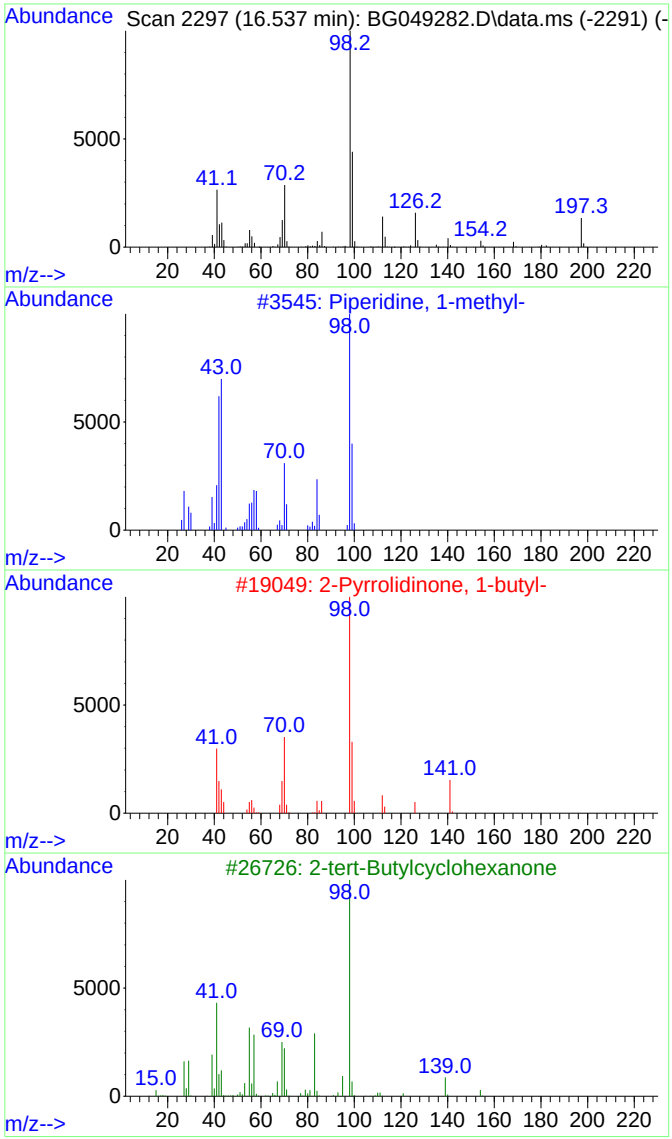
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

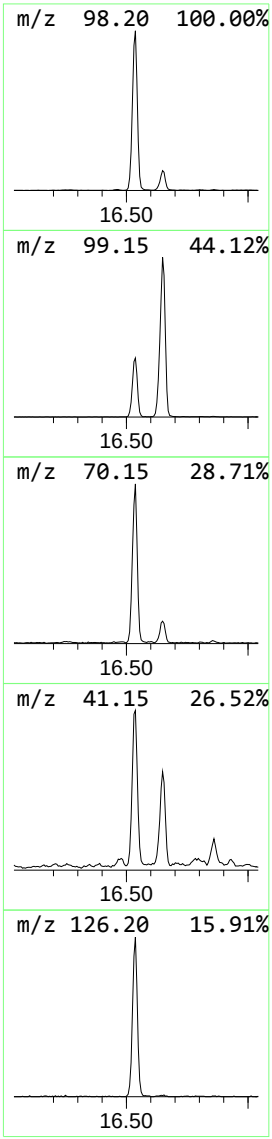
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 Piperidine, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.540	73.17 ng/ul	1807160	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Piperidine, 1-methyl-	99	C6H13N	000626-67-5	53
2	2-Pyrrolidinone, 1-butyl-	141	C8H15NO	003470-98-2	45
3	2-tert-Butylcyclohexanone	154	C10H18O	001728-46-7	43
4	Piperidine, 1-methyl-	99	C6H13N	000626-67-5	42
5	2-Pyrrolidinemethanamine, 1-ethyl-	128	C7H16N2	026116-12-1	38





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ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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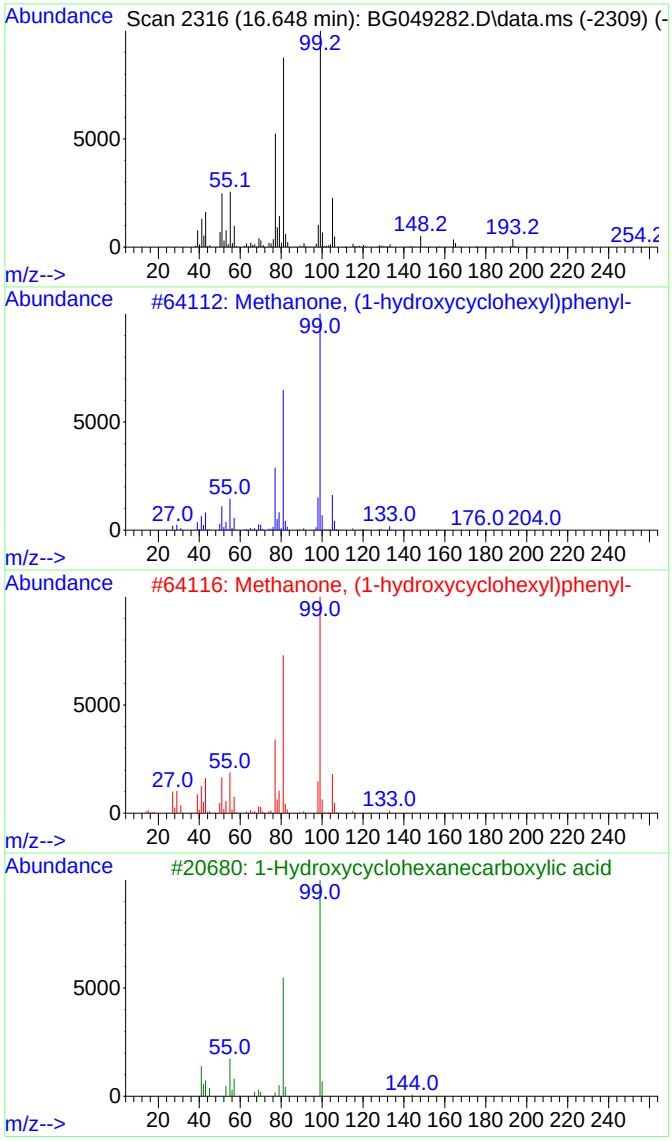
Quant Title : SVOA CALIBRATION

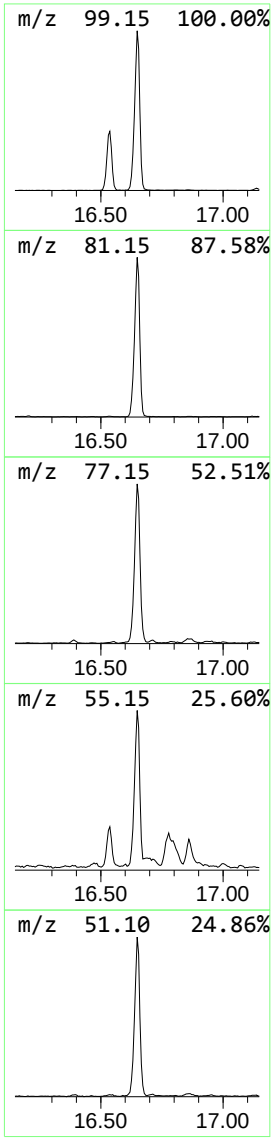
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 Methanone, (1-hydroxycycloh... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.650	153.48 ng/ul	3791010	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	74
2	Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	72
3	1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	47
4	Phosphonofluoridic acid, methyl-...	194	C8H16FO2P	113548-86-0	47
5	Furan, tetrahydro-2,2-dimethyl-5...	156	C10H20O	033978-70-0	47





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Instrument : BNA_G
ClientSampleId : DBK43

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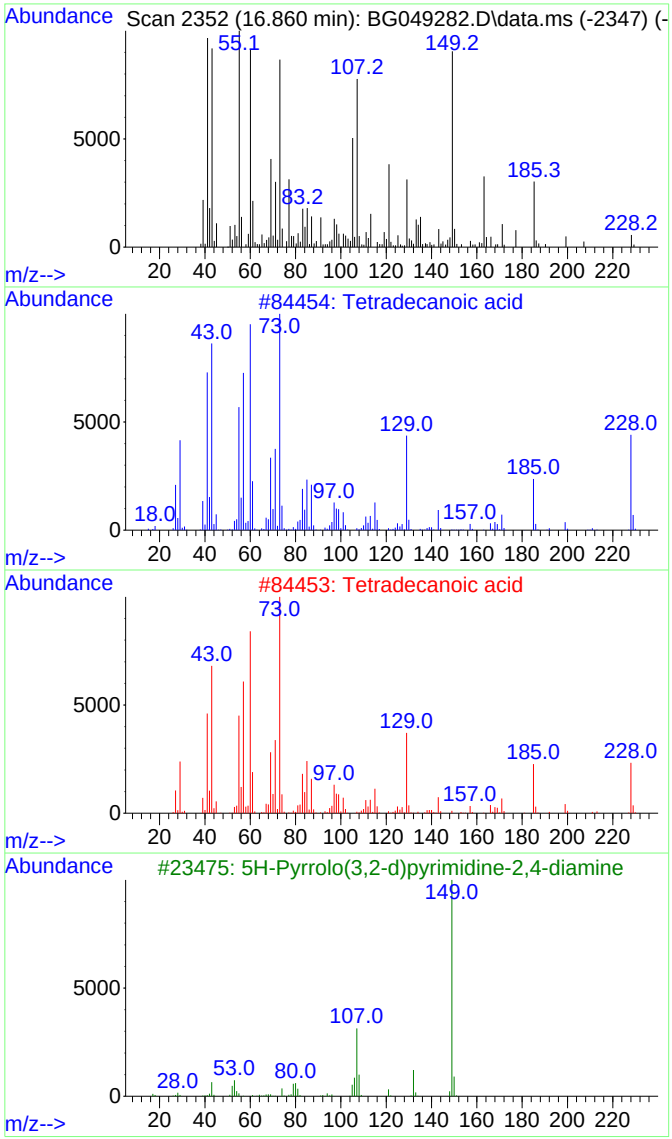
Quant Title : SVOA CALIBRATION

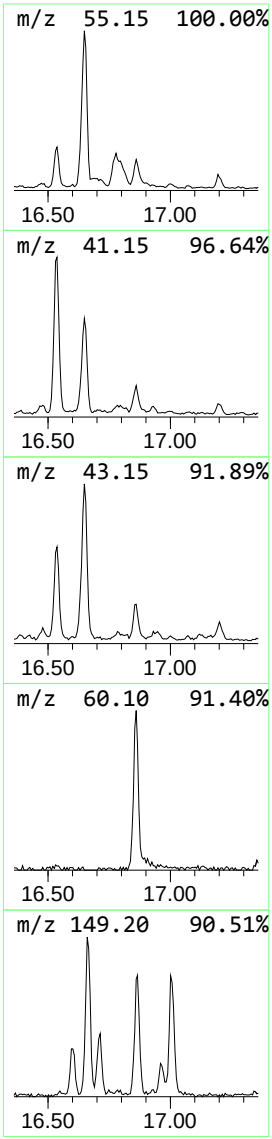
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 36 Tetradecanoic acid Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.860	23.58 ng/ul	582321	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	38
3	5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	38
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	35
5	Tridecanoic acid	214	C13H26O2	000638-53-9	30





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Instrument : BNA_G
ClientSampleId : DBK43

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Quant Title : SVOA CALIBRATION

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

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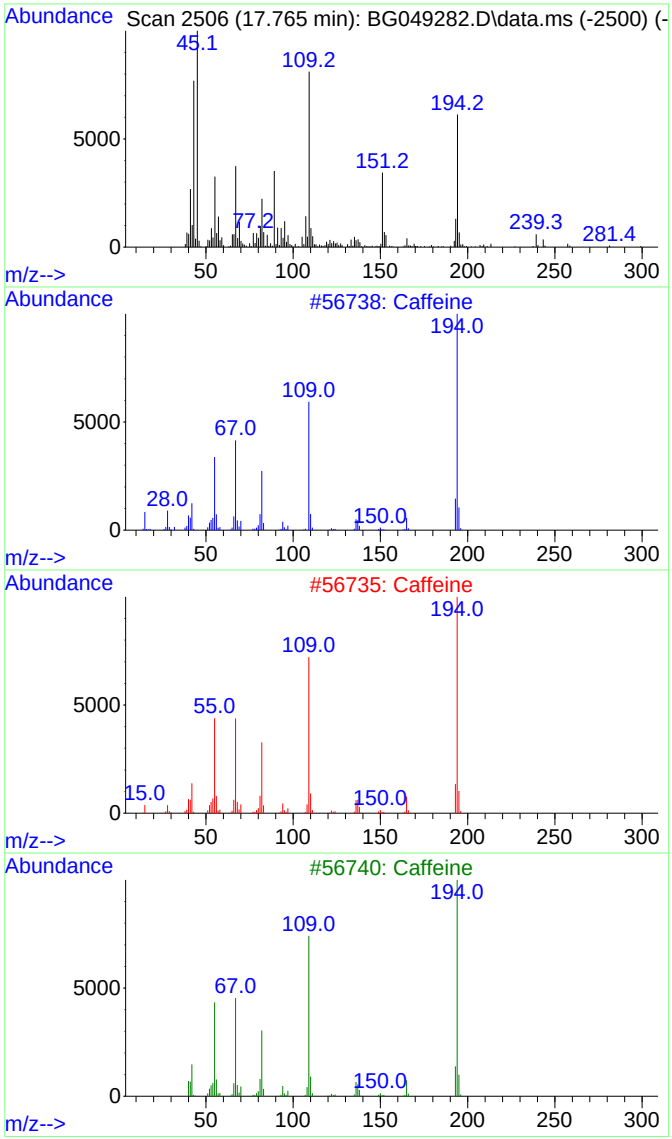
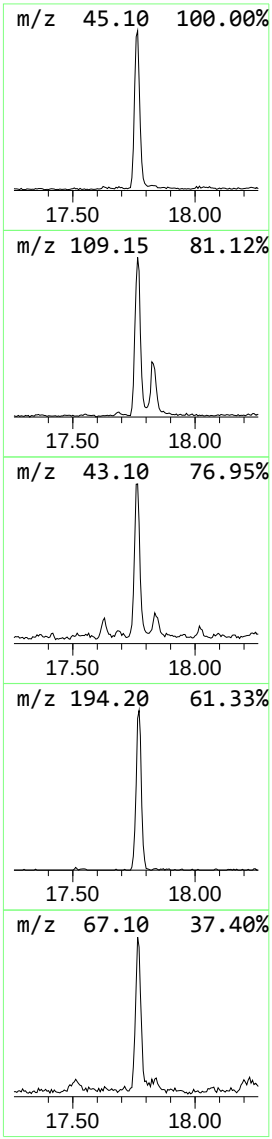
Peak Number 39 Caffeine Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.760	30.69 ng/ul	757976	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Caffeine	194	C8H10N4O2	000058-08-2	78
2	Caffeine	194	C8H10N4O2	000058-08-2	78
3	Caffeine	194	C8H10N4O2	000058-08-2	70
4	Caffeine	194	C8H10N4O2	000058-08-2	50
5	1,4-Dimethyl-4,5,7,8-tetrahydroi...	194	C8H10N4O2	130063-15-9	30





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Instrument : BNA_G
ClientSampleId : DBK43

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Quant Title : SVOA CALIBRATION

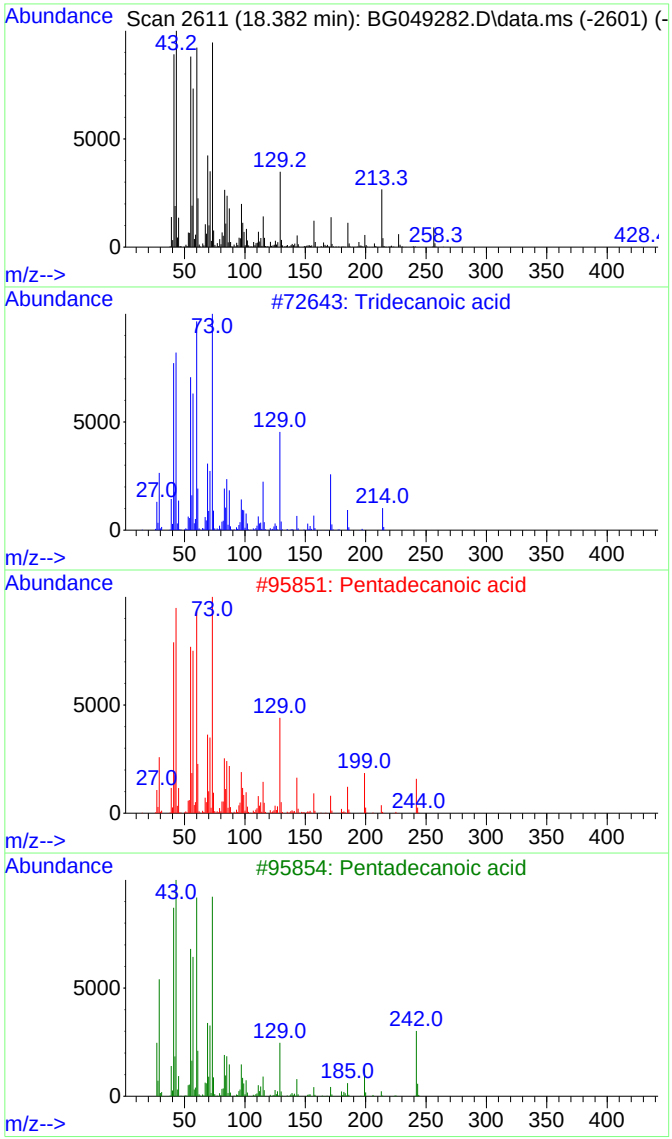
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

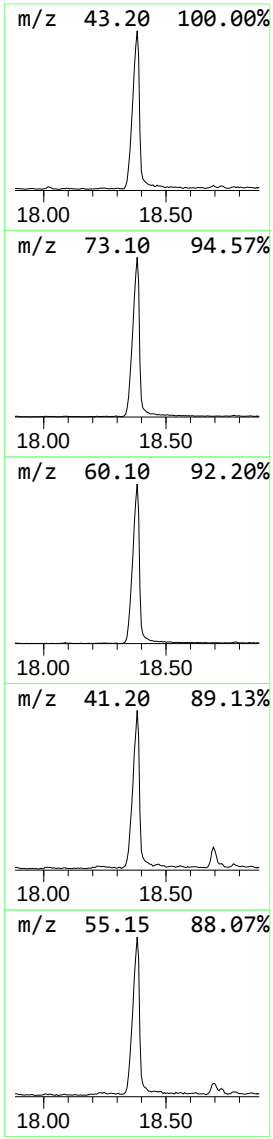
****

Peak Number 41 Tridecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.380	184.53 ng/ul	4557850	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	94
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
5	Tridecanoic acid	214	C13H26O2	000638-53-9	64





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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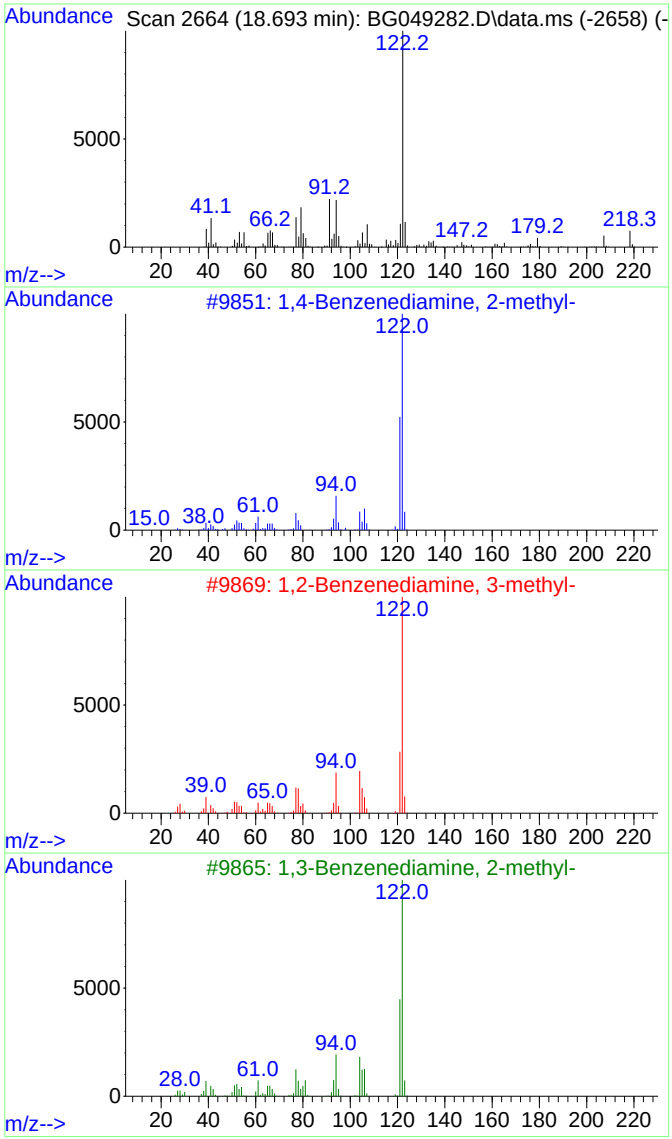
Quant Title : SVOA CALIBRATION

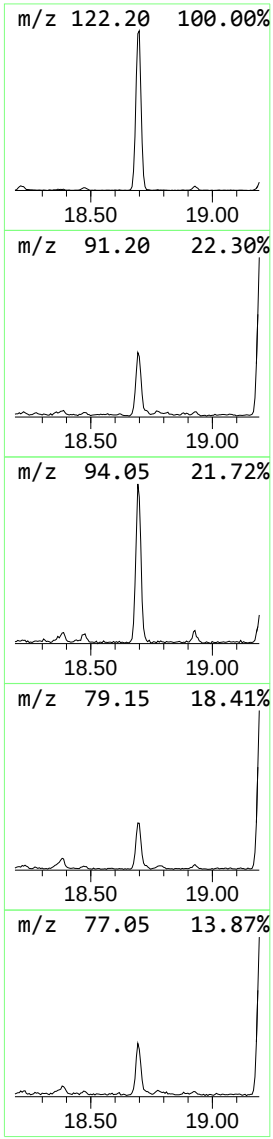
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 42 1,4-Benzenediamine, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.690	44.22 ng/ul	1092270	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenediamine, 2-methyl-	122	C7H10N2	000095-70-5	64
2	1,2-Benzenediamine, 3-methyl-	122	C7H10N2	002687-25-4	64
3	1,3-Benzenediamine, 2-methyl-	122	C7H10N2	000823-40-5	64
4	1,2-Benzenediamine, 4-methyl-	122	C7H10N2	000496-72-0	64
5	1,2-Benzenediamine, 4-methyl-	122	C7H10N2	000496-72-0	64





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Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

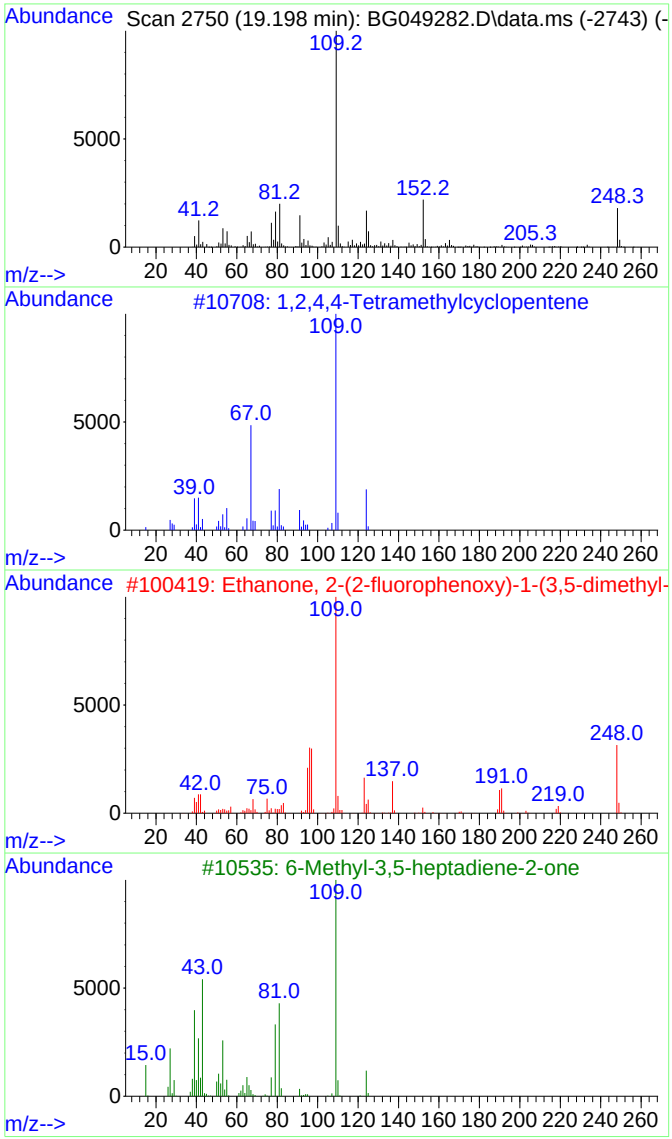
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

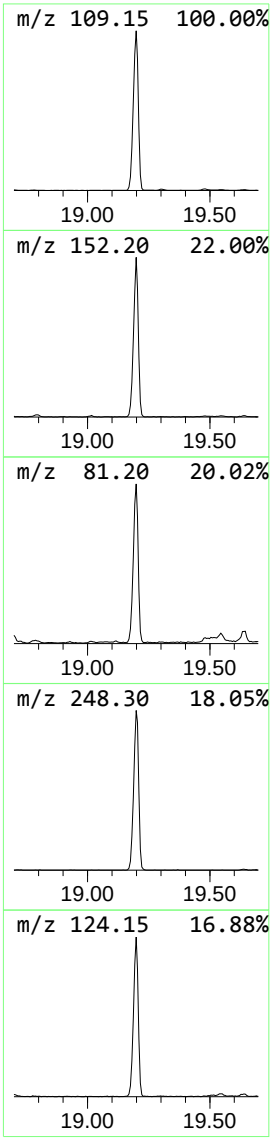
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 44 1,2,4,4-Tetramethylcyclopen... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.200	183.42 ng/ul	4530480	Phenanthrene-d10	17.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,4-Tetramethylcyclopentene	124	C9H16	065378-76-9	70
2	Ethanone, 2-(2-fluorophenoxy)-1-...	248	C13H13FN2O2	344956-93-0	64
3	6-Methyl-3,5-heptadiene-2-one	124	C8H12O	001604-28-0	53
4	2,5-Dimethylhex-5-en-3-yn-2-ol	124	C8H12O	1000302-74-9	53
5	2-Cyclopenten-1-one, 3,4,4-trime...	124	C8H12O	030434-65-2	53





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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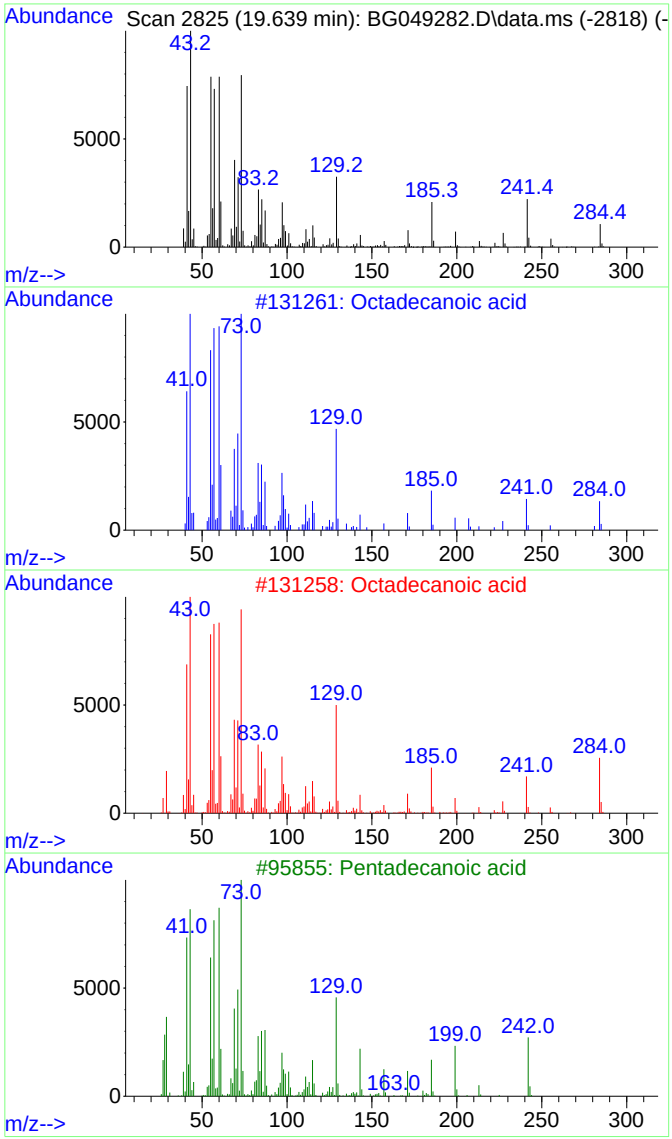
Quant Title : SVOA CALIBRATION

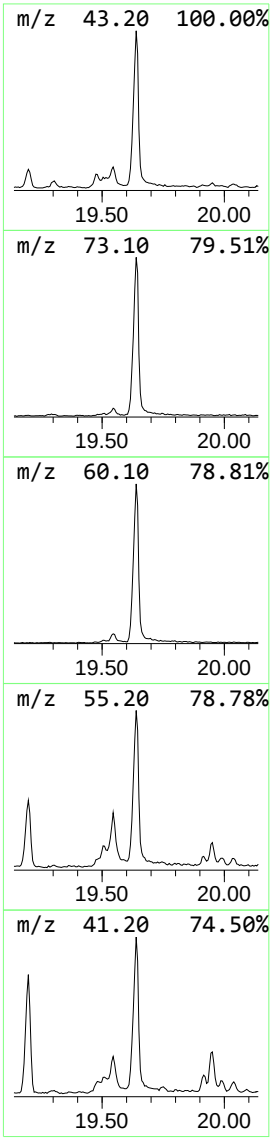
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 48 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.640	121.84 ng/ul	2835370	Chrysene-d12	21.754

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	87
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80
5	Octadecanoic acid	284	C18H36O2	000057-11-4	74





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Instrument : BNA_G
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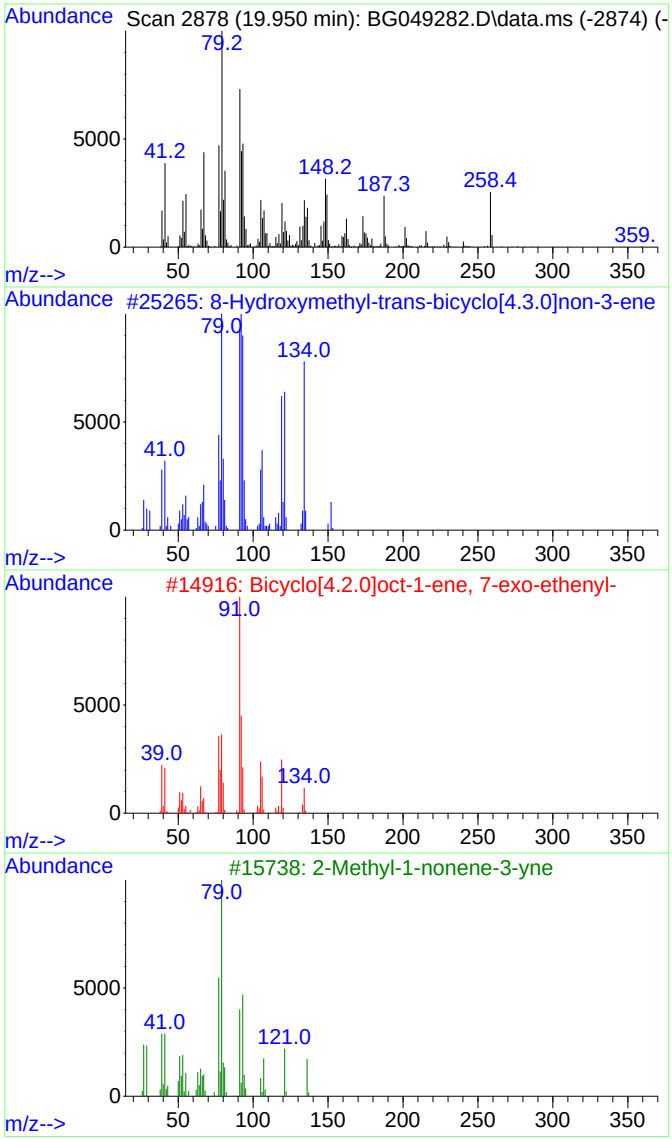
Quant Title : SVOA CALIBRATION

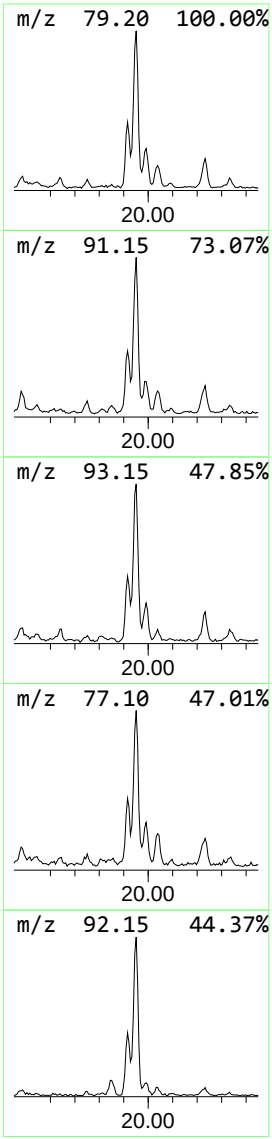
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 51 8-Hydroxymethyl-trans-bicyc... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.950	55.70 ng/ul	1296330	Chrysene-d12	21.754

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Hydroxymethyl-trans-bicyclo[4....	152	C10H16O	1000099-21-8	53
2	Bicyclo[4.2.0]oct-1-ene, 7-exo-e...	134	C10H14	1000142-18-2	43
3	2-Methyl-1-nonene-3-yne	136	C10H16	070058-00-3	38
4	Cyclohexane, 1-butenylidene-	136	C10H16	036144-40-8	38
5	2-Methyl-1-nonene-3-yne	136	C10H16	070058-00-3	38





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Quant Title : SVOA CALIBRATION

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

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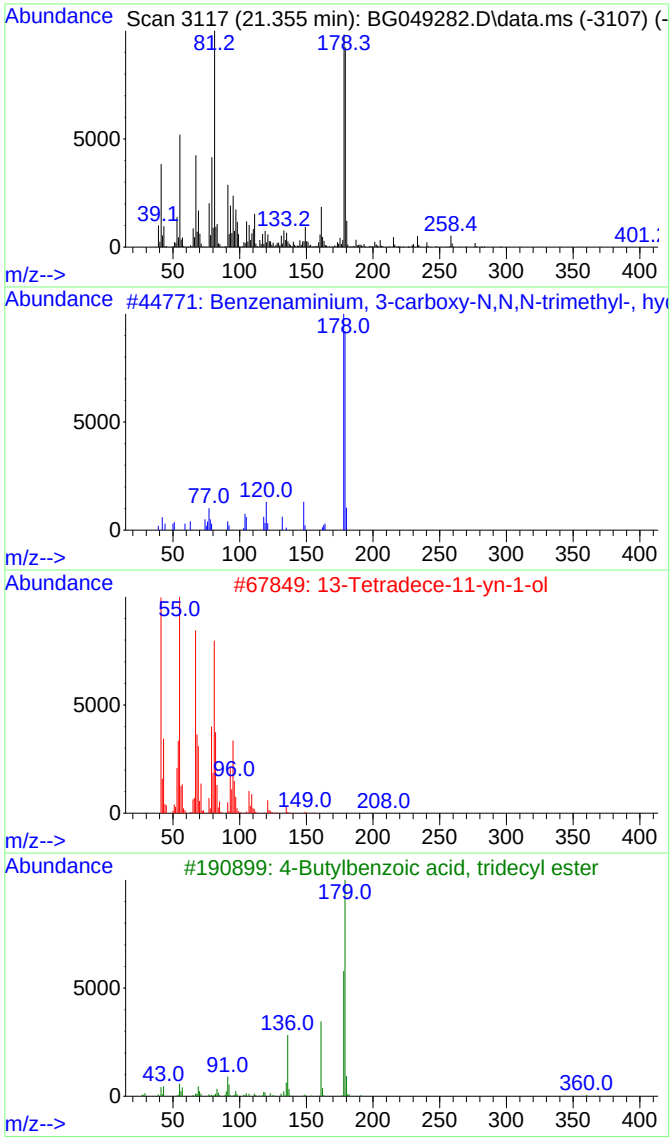
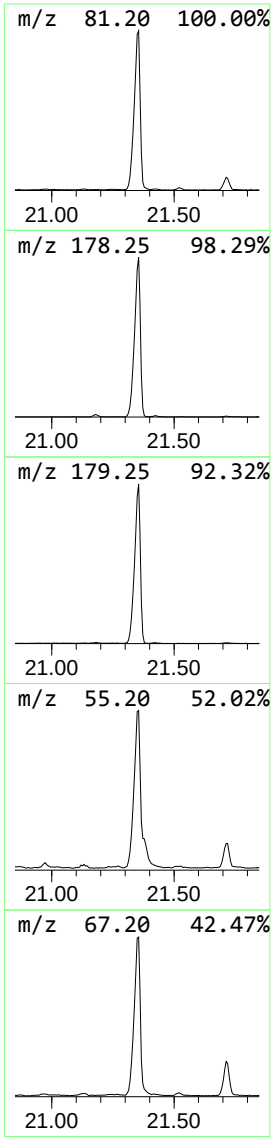
Peak Number 60 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.350	378.86 ng/ul	8816540	Chrysene-d12	21.754

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenaminium, 3-carboxy-N,N,N-t...	179	C10H13NO2	033192-03-9	35
2	13-Tetradece-11-yn-1-ol	208	C14H24O	1000131-00-4	25
3	4-Butylbenzoic acid, tridecyl ester	360	C24H40O2	1000292-32-6	22
4	2,4,6-Trimethyl-7-oxo-4,7-dihydr...	179	C7H9N5O	025623-78-3	22
5	5-Nitrobenzofuran-2-one	179	C8H5NO4	021997-23-9	22





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Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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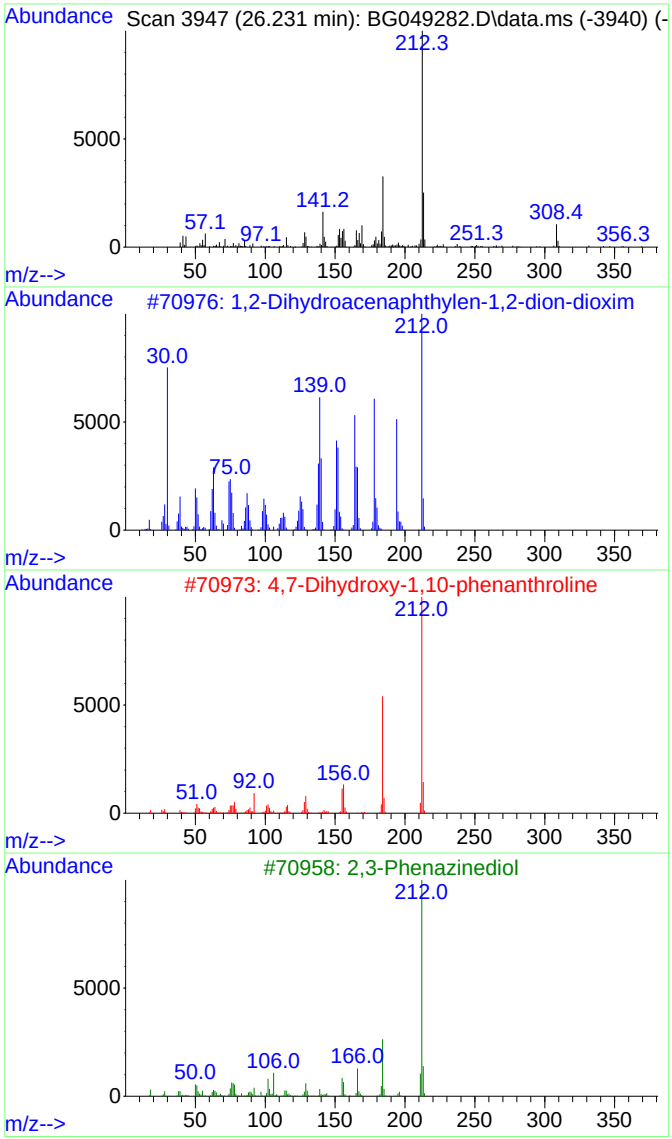
Quant Title : SVOA CALIBRATION

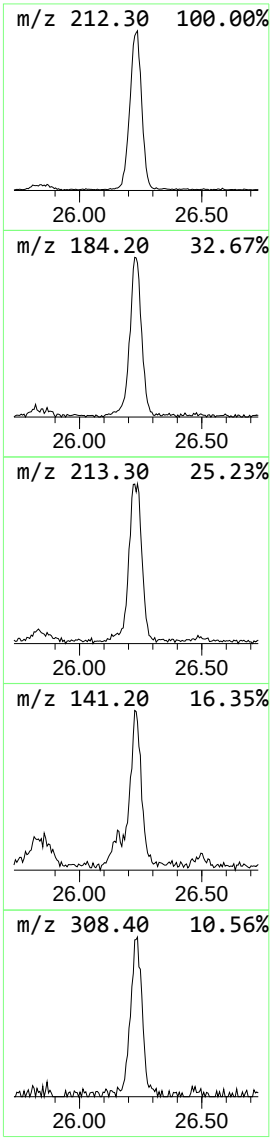
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 72 1,2-Dihydroacenaphthylen-1,... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.230	32.27 ng/ul	914989	Perylene-d12	25.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Dihydroacenaphthylen-1,2-dio...	212	C12H8N2O2	001932-08-7	60
2	4,7-Dihydroxy-1,10-phenanthroline	212	C12H8N2O2	003922-40-5	53
3	2,3-Phenazinediol	212	C12H8N2O2	019220-18-9	53
4	1,3-Benzodioxole-5-carbonitrile,...	212	C12H8N2O2	1000351-67-6	47
5	1-(4-Methoxyphenyl)-2-(4-pyridaz...	212	C13H12N2O	1000279-82-6	47





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Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

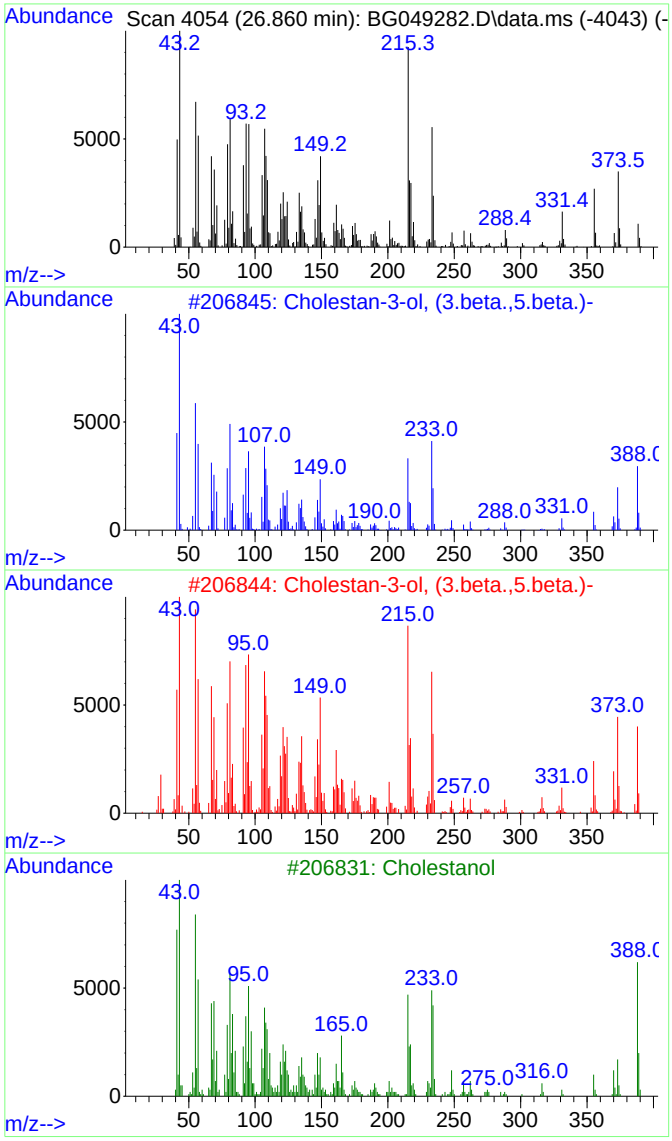
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

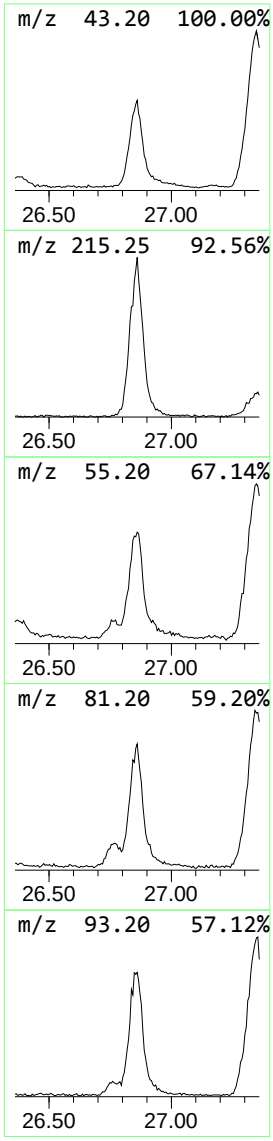
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 74 Cholestan-3-ol, (3.beta.,5.... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.860	135.21 ng/ul	3834380	Perylene-d12	25.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	74
2	Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	64
3	Cholestanol	388	C27H48O	000080-97-7	60
4	Cholestanol	388	C27H48O	000080-97-7	49
5	2-((1S,3S)-3-hydroxycyclohexyl)-...	332	C22H36O2	1000372-26-3	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

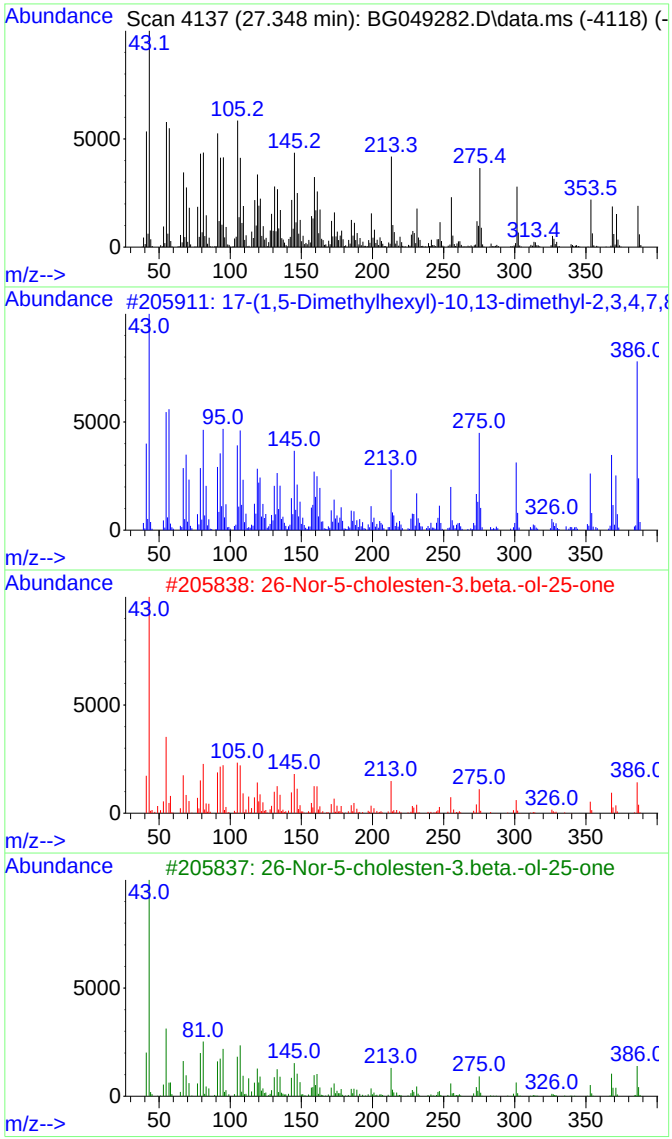
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

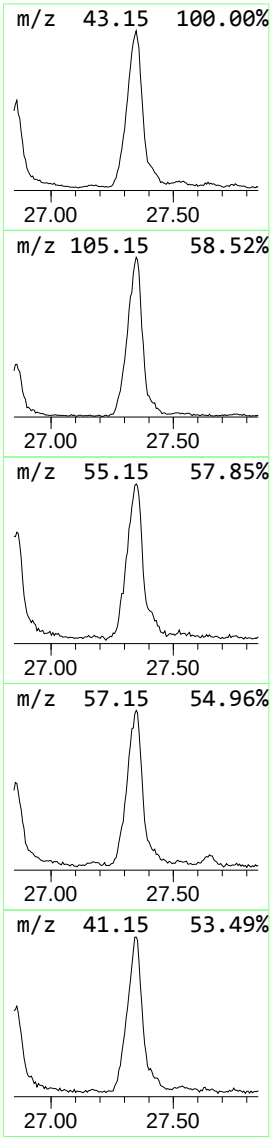
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 75 17-(1,5-Dimethylhexyl)-10,1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.350	265.50 ng/ul	7529080	Perylene-d12	25.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-38-4	99
2	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
3	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	98
4	Cholesterol	386	C27H46O	000057-88-5	98
5	Cholest-5-en-3-ol, (3.alpha.)-	386	C27H46O	000474-77-1	62





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Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

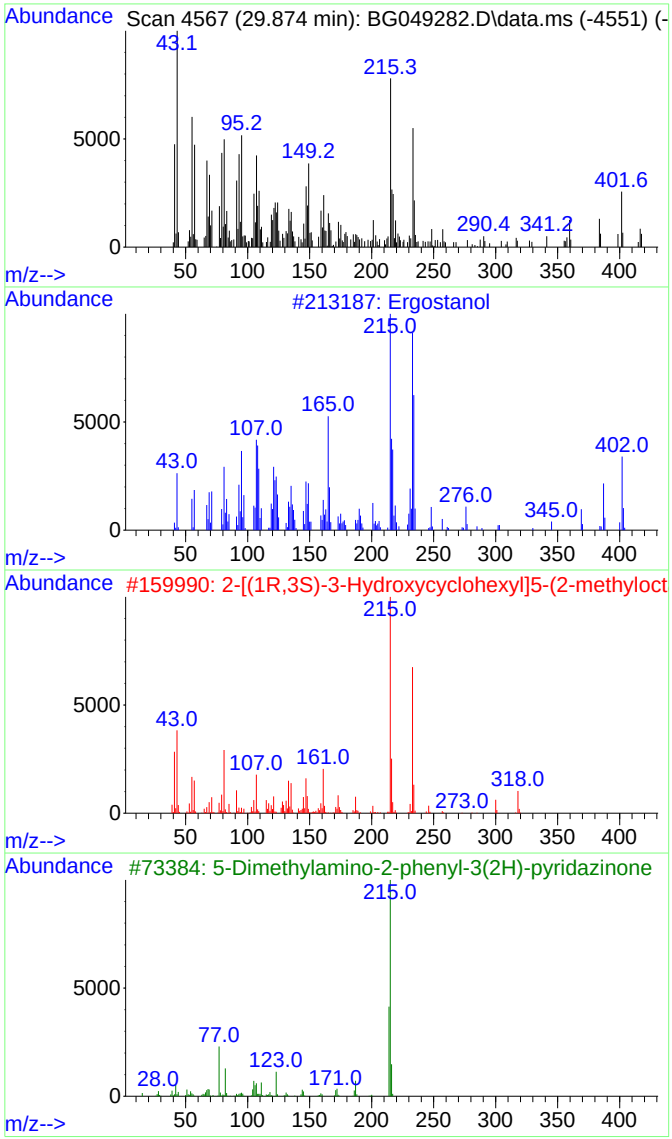
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

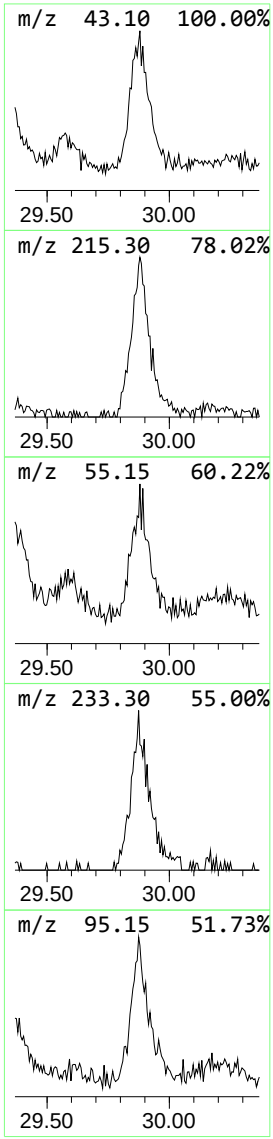
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 79 Ergostanol Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.870	26.61 ng/ul	754576	Perylene-d12	25.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ergostanol	402	C28H50O	006538-02-9	53
2	2-[(1R,3S)-3-Hydroxycyclohexyl]5...	318	C21H34O2	070434-82-1	38
3	5-Dimethylamino-2-phenyl-3(2H)-p...	215	C12H13N3O	1000244-25-4	20
4	Quinoxaline, 2-(2,4-dimethylphen...	234	C16H14N2	060814-26-8	15
5	5-Amino-2-(2-methylpropyl)-1-pen...	285	C16H23N5	1000326-96-2	14





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Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

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

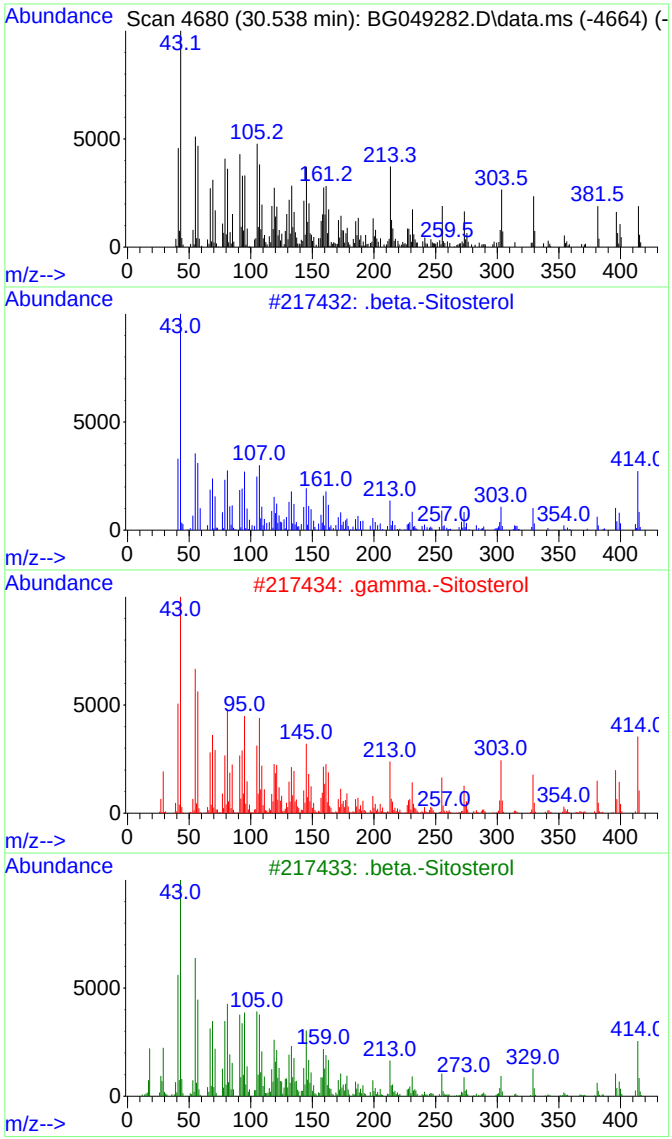
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

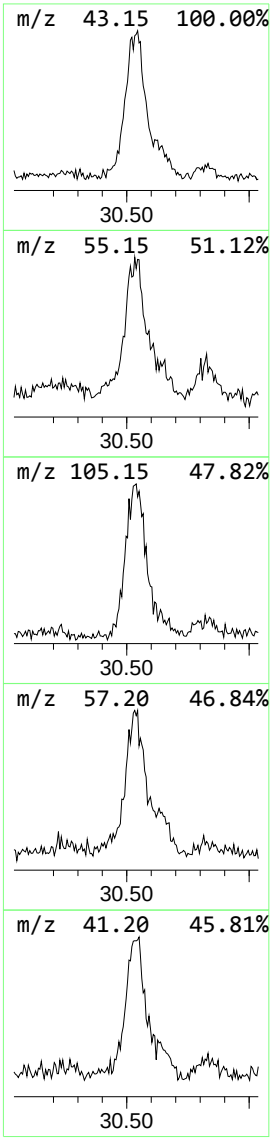
****

Peak Number 80 .beta.-Sitosterol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.540	59.45 ng/ul	1685800	Perylene-d12	25.044

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Sitosterol	414	C29H50O	000083-46-5	93
2	.gamma.-Sitosterol	414	C29H50O	000083-47-6	89
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	49
4	.gamma.-Sitosterol	414	C29H50O	000083-47-6	47
5	Ergost-7-en-3-ol	400	C28H48O	116179-22-7	35





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
Data File : BG049282.D
Acq On : 11 Jul 2021 2:38
Operator : CG/JU
Sample : M2914-14
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBK43

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.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
(DEL) Alkane: C...	3.120	20.3	ng/ul	232123	1	8.076	228365	20.0
Butanoic acid, ...	5.000	28.7	ng/ul	328122	1	8.076	228365	20.0
3-Hexanone, 2-m...	5.900	52.6	ng/ul	601064	1	8.076	228365	20.0
Cyclohexanol	5.950	102.7	ng/ul	1172410	1	8.076	228365	20.0
Cyclohexanone	6.100	35.4	ng/ul	404380	1	8.076	228365	20.0
Benzene, 1,2,3-...	7.720	36.0	ng/ul	411025	1	8.076	228365	20.0
unknown-01	7.850	79.4	ng/ul	907087	1	8.076	228365	20.0
unknown-02	10.120	23.7	ng/ul	567854	2	10.896	479247	20.0
1-(1-Methoxypro...	10.360	32.9	ng/ul	788866	2	10.896	479247	20.0
Benzoic acid	10.710	73.0	ng/ul	1750350	2	10.896	479247	20.0
1-Phenoxypropan...	11.690	741.6	ng/ul	17770200	2	10.896	479247	20.0
Benzeneacetic acid	11.870	32.5	ng/ul	778120	2	10.896	479247	20.0
(DEL) Alkane: S...	12.300	7.1	ng/ul	171122	2	10.896	479247	20.0
Indole	12.510	31.6	ng/ul	758388	2	10.896	479247	20.0
Cyclohexanone, ...	15.120	497.6	ng/ul	12097300	3	14.721	486180	20.0
2H-Benzocyclohe...	15.330	39.0	ng/ul	948627	3	14.721	486180	20.0
N-Cyclohexyl-2-...	15.590	2797.7	ng/ul	68008700	3	14.721	486180	20.0
Piperidine, 1-m...	16.540	73.2	ng/ul	1807160	4	17.471	493996	20.0
Methanone, (1-h...	16.650	153.5	ng/ul	3791010	4	17.471	493996	20.0
Tetradecanoic acid	16.860	23.6	ng/ul	582321	4	17.471	493996	20.0
Caffeine	17.760	30.7	ng/ul	757976	4	17.471	493996	20.0
Tridecanoic acid	18.380	184.5	ng/ul	4557850	4	17.471	493996	20.0
1,4-Benzenediam...	18.690	44.2	ng/ul	1092270	4	17.471	493996	20.0
1,2,4,4-Tetrame...	19.200	183.4	ng/ul	4530480	4	17.471	493996	20.0
Octadecanoic acid	19.640	121.8	ng/ul	2835370	5	21.754	465429	20.0
8-Hydroxymethyl...	19.950	55.7	ng/ul	1296330	5	21.754	465429	20.0
unknown-03	21.350	378.9	ng/ul	8816540	5	21.754	465429	20.0
1,2-Dihydroacen...	26.230	32.3	ng/ul	914989	6	25.044	567166	20.0
Cholestan-3-ol,...	26.860	135.2	ng/ul	3834380	6	25.044	567166	20.0
17-(1,5-Dimethy...	27.350	265.5	ng/ul	7529080	6	25.044	567166	20.0
Ergostanol	29.870	26.6	ng/ul	754576	6	25.044	567166	20.0
.beta.-Sitosterol	30.540	59.5	ng/ul	1685800	6	25.044	567166	20.0