LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\
Data File : BG041946.D
Acq On : 4 Aug 2019 17:36
Operator : HP/JU
Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

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\SOM-EPA-BG080319MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.159	7	12	38	rVB	1141370	2044079	100.00%	9.852%
2	3.417	54	56	64	rVB5	11702	20102	0.98%	0.097%
3	4.093	166	171	176	rVB3	7500	12395	0.61%	0.060%
4	4.187	182	187	193	rVB4	2913	4920	0.24%	0.024%
5	5.109	340	344	349	rVB4	2514	3783	0.19%	0.018%
6	6.096	508	512	514	rBV4	1622	2124	0.10%	0.010%
7	6.672	606	610	612	rBV3	1728	2045	0.10%	0.010%
8	7.183	690	697	714	rBV	212252	423819	20.73%	2.043%
9	7.354	720	726	739	rBV	153935	285441	13.96%	1.376%
10	7.565	755	762	776	rBV	227531	432755	21.17%	2.086%
11	8.035	835	842	851	rBV	180785	321581	15.73%	1.550%
12	8.100	851	853	857	rVB3	3102	4054	0.20%	0.020%
13	8.740	955	962	973	rBV	247212	444843	21.76%	2.144%
14	8.864	977	983	988	rVV5	2631	5273	0.26%	0.025%
15	9.116	1020	1026	1030	rBV4	1817	3319	0.16%	0.016%
16	9.199	1032	1040	1050	rBV	220878	412662	20.19%	1.989%
17	9.369	1067	1069	1072	rVB3	2197	2379	0.12%	0.011%
18	9.933	1155	1165	1179	rBV	196915	370698	18.14%	1.787%
19	10.474	1249	1257	1274	rBV	299230	601260	29.41%	2.898%
20	10.861	1312	1323	1335	rBV	261147	471950	23.09%	2.275%
21	10.991	1337	1345	1357	rVV	223083	442693	21.66%	2.134%
22	12.454	1589	1594	1599	rVB2	5244	8679	0.42%	0.042%
23	13.834	1822	1829	1833	rBV2	4370	9230	0.45%	0.044%
24	14.093	1867	1873	1877	rBV	552423	832578	40.73%	4.013%
25	14.140	1877	1881	1897	rVB	264571	423437	20.72%	2.041%
26	14.387	1914	1923	1935	rBV	662216	1074969	52.59%	5.181%
27	14.698	1968	1976	1984	rBV	387535	665222	32.54%	3.206%
28	14.874	1999	2006	2021	rBV	163069	315761	15.45%	1.522%
29	15.098	2042	2044	2051	rVB6	3541	4674	0.23%	0.023%
30	15.497	2107	2112	2113	rBV2	1906	2431	0.12%	0.012%
31	15.550	2116	2121	2125	rVB4	2427	3686	0.18%	0.018%
32	15.691	2138	2145	2158	rBV	873716	1372232	67.13%	6.614%
33	15.803	2158	2164	2175	rVV	179618	287951	14.09%	1.388%
34	15.932	2181	2186	2193	rVB	10815	17129	0.84%	0.083%

LSC Area Percent Report

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Operator : HP/JU
Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

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\SOM-EPA-BG080319MA.M
Title : SVOA CALIBRATION

35	16.337	2251	2255	2258	rBV3	1790	2610	0.13%	0.013%
36	16.384	2258	2263	2266	rBV3	3179	6278	0.31%	0.030%
37	16.473	2275	2278	2285	rVB6	5005	6982	0.34%	0.034%
38	16.849	2339	2342	2347	rBV4	2993	5458	0.27%	0.026%
39	17.148	2388	2393	2396	rVV7	6106	8517	0.42%	0.041%
40	17.207	2399	2403	2404	rBV2	3666	3072	0.15%	0.015%
41	17.348	2423	2427	2430	rBV2	3677	5785	0.28%	0.028%
42	17.389	2432	2434	2435	rBV2	2598	2468	0.12%	0.012%
43	17.448	2435	2444	2450	rBV	517586	813856	39.82%	3.923%
44	17.489	2450	2451	2455	rVV	20011	17358	0.85%	0.084%
45	17.548	2455	2461	2469	rVB	948976	1455891	71.22%	7.017%
46	17.830	2503	2509	2512	rBV8	10329	18299	0.90%	0.088%
47	17.941	2526	2528	2531	rBV3	2956	3220	0.16%	0.016%
48	18.035	2542	2544	2548	rVB5	2816	3923	0.19%	0.019%
49	18.071	2548	2550	2551	rBV2	3067	2350	0.11%	0.011%
50	18.112	2555	2557	2558	rBV2	2914	2547	0.12%	0.012%
51	18.176	2566	2568	2571	rBV4	2576	3439	0.17%	0.017%
52	18.206	2571	2573	2575	rBV3	3842	3777	0.18%	0.018%
53	18.288	2583	2587	2588	rBV4	4522	6087	0.30%	0.029%
54	18.341	2590	2596	2611	rVV4	120682	232726	11.39%	1.122%
55	18.482	2619	2620	2622	rBV2	2771	2137	0.10%	0.010%
56	18.517	2622	2626	2631	rVV8	7961	14329	0.70%	0.069%
57	18.599	2638	2640	2642	rBV3	3201	2626	0.13%	0.013%
58	18.635	2643	2646	2648	rBV3	5971	7532	0.37%	0.036%
59	18.699	2655	2657	2658	rBV2	4773	2640	0.13%	0.013%
60	18.793	2671	2673	2675	rVB3	5172	2971	0.15%	0.014%
61	18.817	2675	2677	2678	rBV2	5270	4433	0.22%	0.021%
62	19.263	2751	2753	2757	rVB5	10513	10747	0.53%	0.052%
63	19.463	2784	2787	2790	rBV5	30083	44403	2.17%	0.214%
64	19.498	2790	2793	2797	rVB	45168	57745	2.82%	0.278%
65	19.610	2809	2812	2820	rBV10	28392	55437	2.71%	0.267%
66	19.833	2844	2850	2861	rBV	1126283	1806641	88.38%	8.707%
67	20.092	2891	2894	2898	rVB2	25170	35088	1.72%	0.169%
68	20.503	2960	2964	2976	rVB4	77071	158146	7.74%	0.762%
69	21.038	3051	3055	3059	rVB	75380	93940	4.60%	0.453%
70	21.373	3108	3112	3121	rBV2	98483	188226	9.21%	0.907%
71	21.625	3151	3155	3160	rVB	96378	139957	6.85%	0.675%

LSC Area Percent Report

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Instrument : BNA_G
ClientSampleId : BEP08

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SVOA CALIBRATION

72	21.731	3167	3173	3179	rBV2	518995	876808	42.90%	4.226%
73	22.254	3258	3262	3271	rVB2	76142	140165	6.86%	0.676%
74	22.366	3277	3281	3289	rVB2	92726	168334	8.24%	0.811%
75	23.276	3432	3436	3444	rVB	82870	148158	7.25%	0.714%
76	23.464	3463	3468	3476	rVB	160273	323536	15.83%	1.559%
77	24.099	3572	3576	3585	rVB2	85161	186157	9.11%	0.897%
78	24.780	3684	3692	3708	rVB	555288	1581309	77.36%	7.621%
79	25.010	3723	3731	3739	rBV	279468	761892	37.27%	3.672%

Sum of corrected areas: 20748154

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Operator : HP/JU
Sample : K3962-15
Misc :
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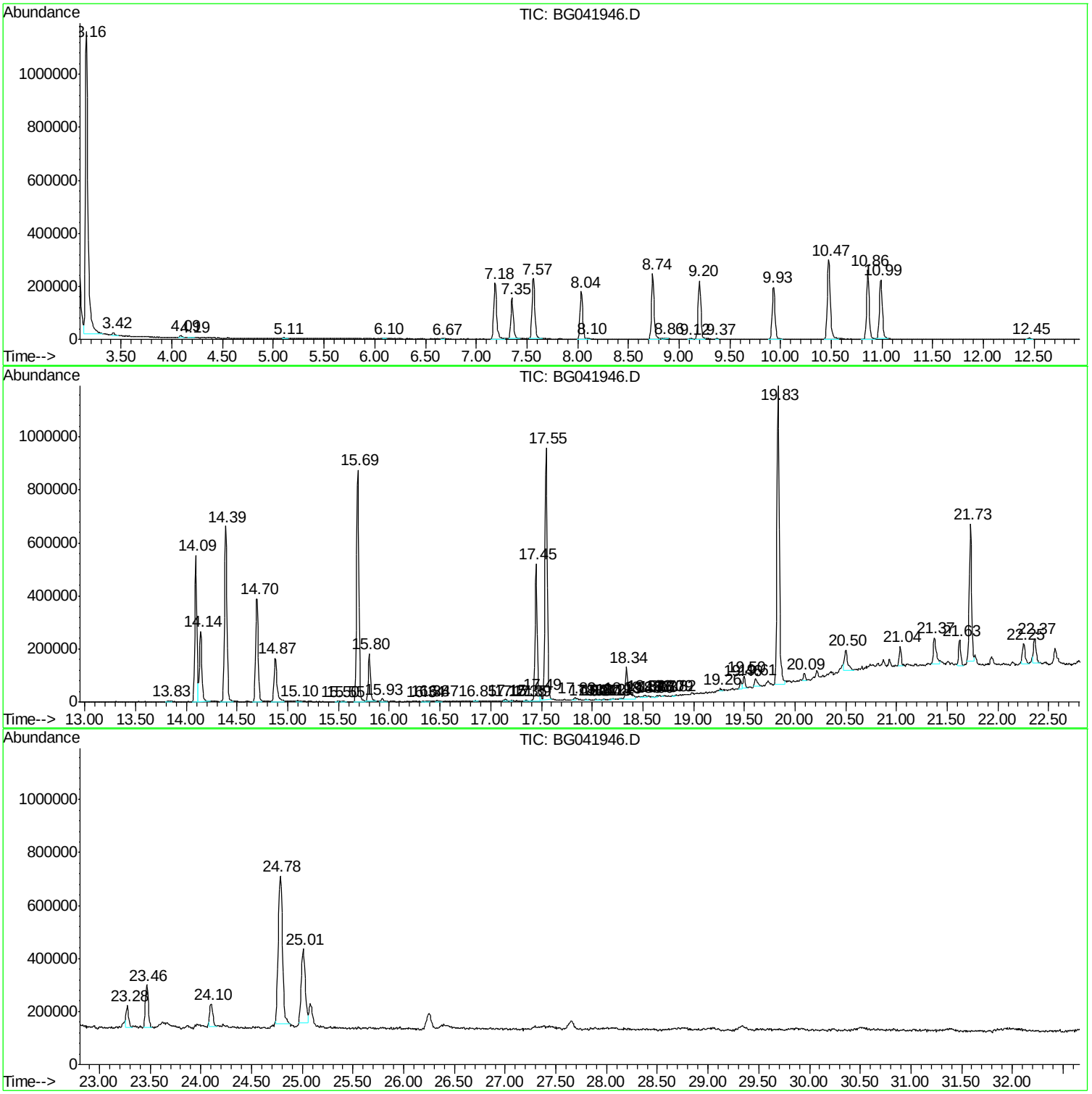
Instrument :
BNA_G
ClientSampleId :
BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\
Data File : BG041946.D
Acq On : 4 Aug 2019 17:36
Operator : HP/JU
Sample : K3962-15
Misc :
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Instrument : BNA_G
ClientSampleId : BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

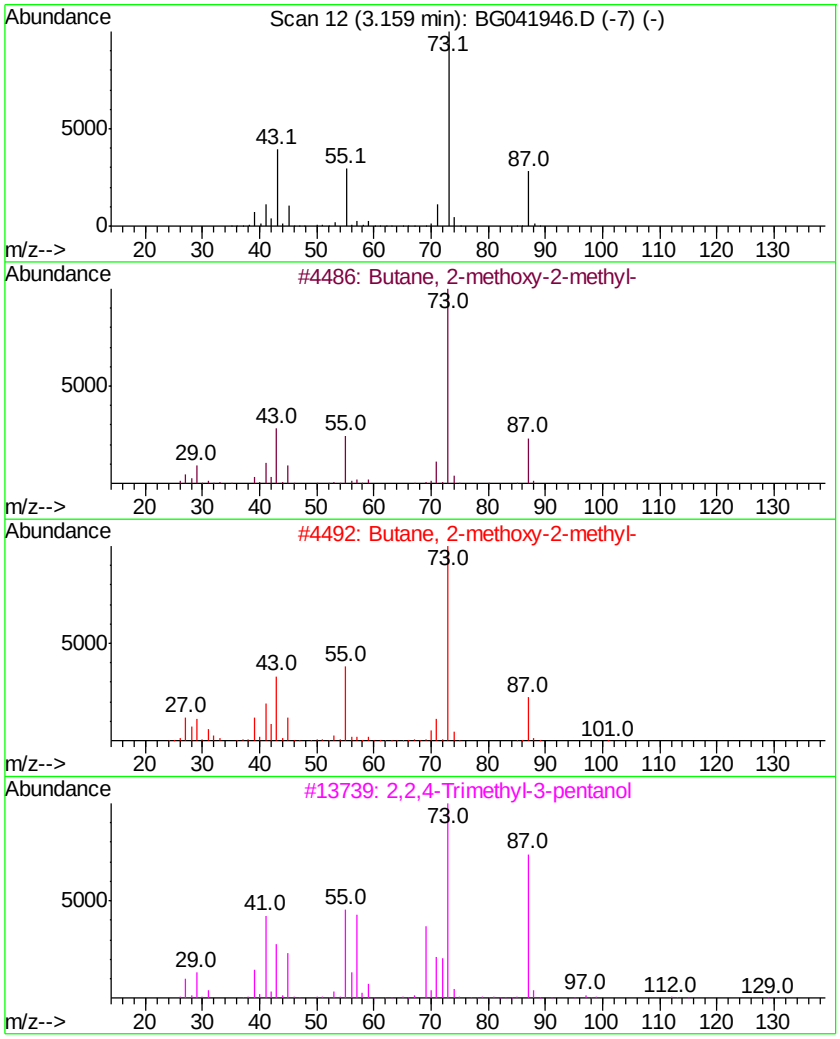
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

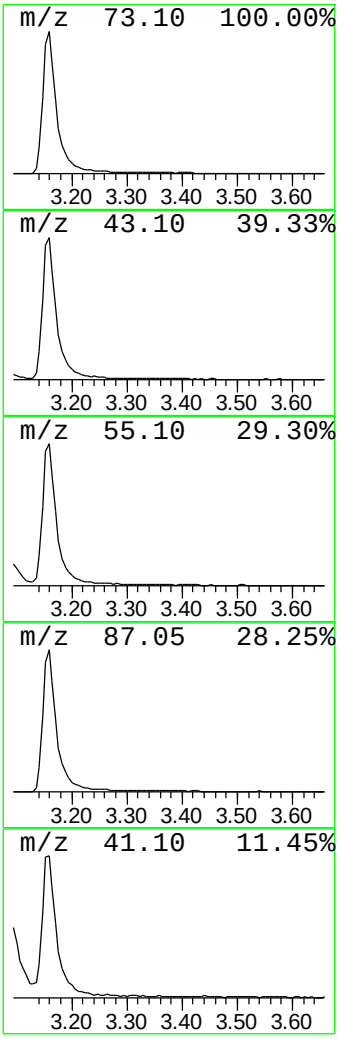
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	127.13 ng/ul	2044080	1,4-Dichlorobenzene-d4	8.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	53
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	33
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Library Search Compound Report

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Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

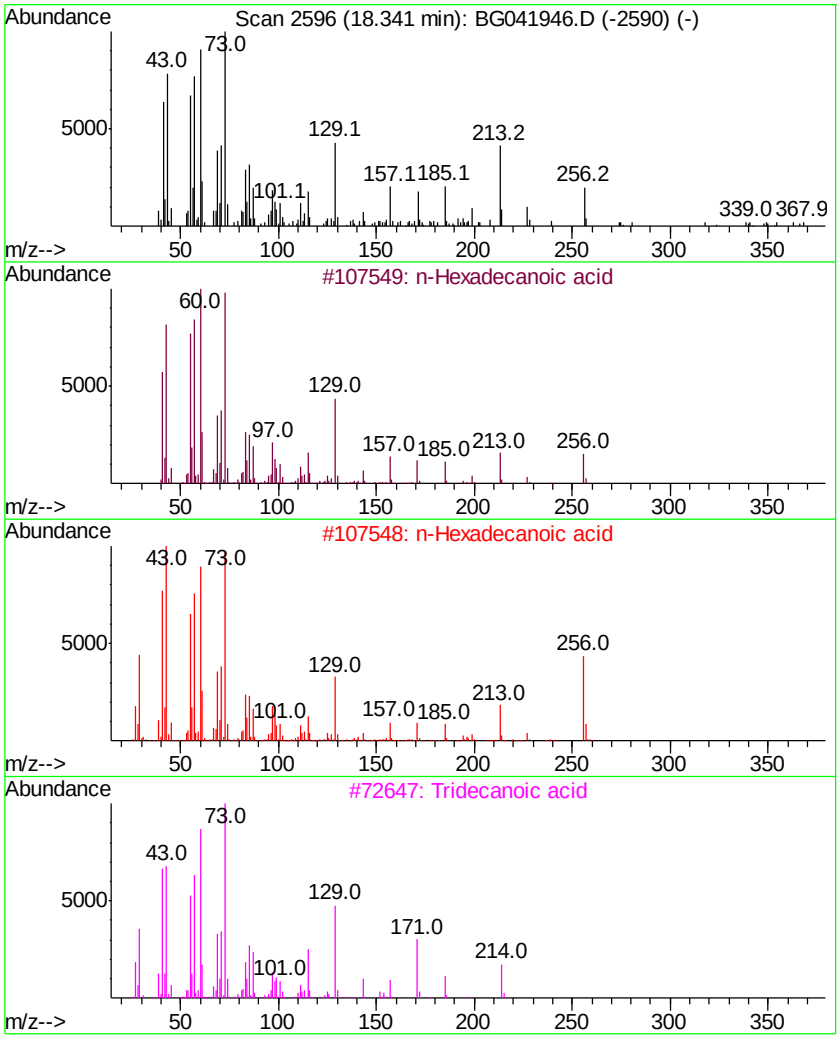
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

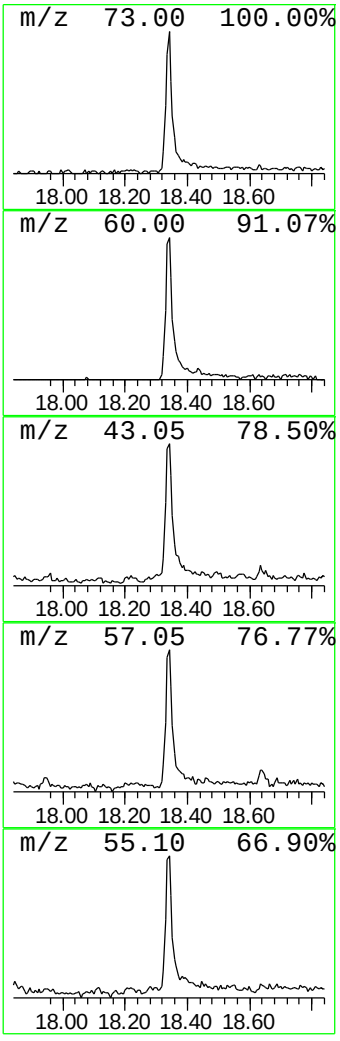
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	5.72 ng/ul	232726	Phenanthrene-d10	17.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	Tridecanoic acid	214	C13H26O2	000638-53-9	93
4	Tridecanoic acid	214	C13H26O2	000638-53-9	91
5	Tridecanoic acid	214	C13H26O2	000638-53-9	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\
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ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

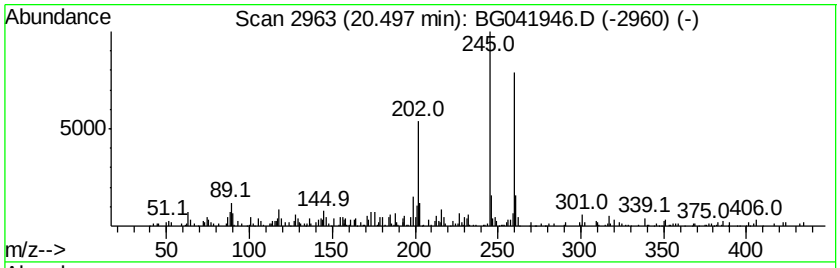
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

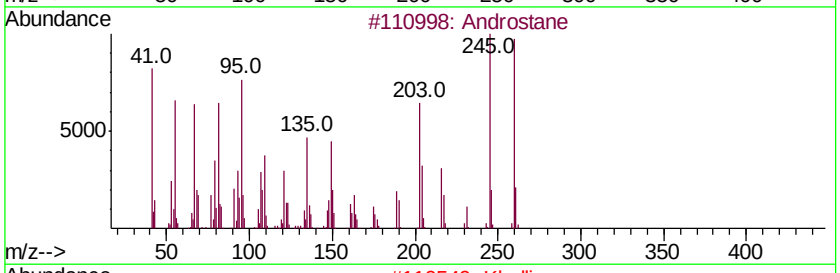
****

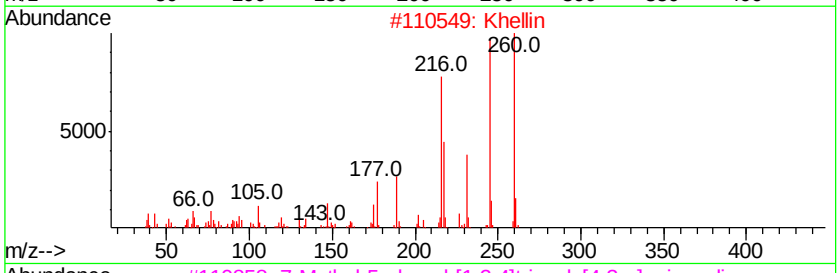
Peak Number 3 Androstane Concentration Rank 7

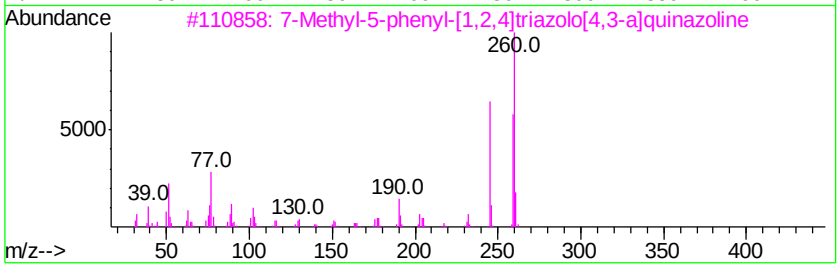
R.T.	EstConc	Area	Relative to ISTD	R.T.
20.50	3.61 ng/ul	158146	Chrysene-d12	21.73

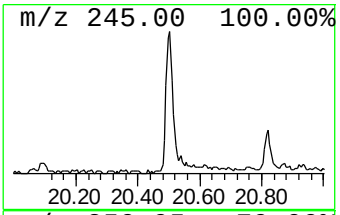
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Androstane	260	C19H32	024887-75-0	98
2	Khellin	260	C14H12O5	000082-02-0	64
3	7-Methyl-5-phenyl-[1,2,4]triazol...	260	C16H12N4	1000318-31-3	46
4	6-[1-[4-Fluorophenyl]ethyl]-1,3-...	260	C15H13FO3	1000211-52-8	43
5	Benzo[a]anthracene, 6,7-dimethyl...	260	C20H20	063020-36-0	38

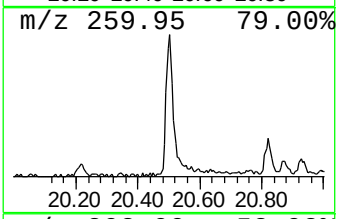


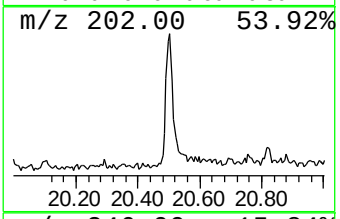


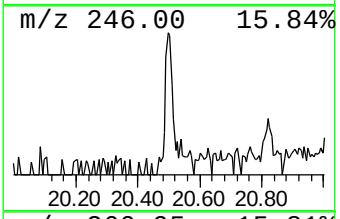


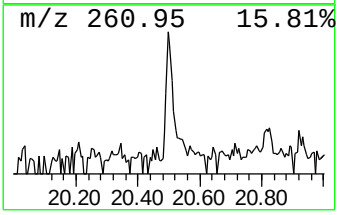












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Quant Title : SVOA CALIBRATION

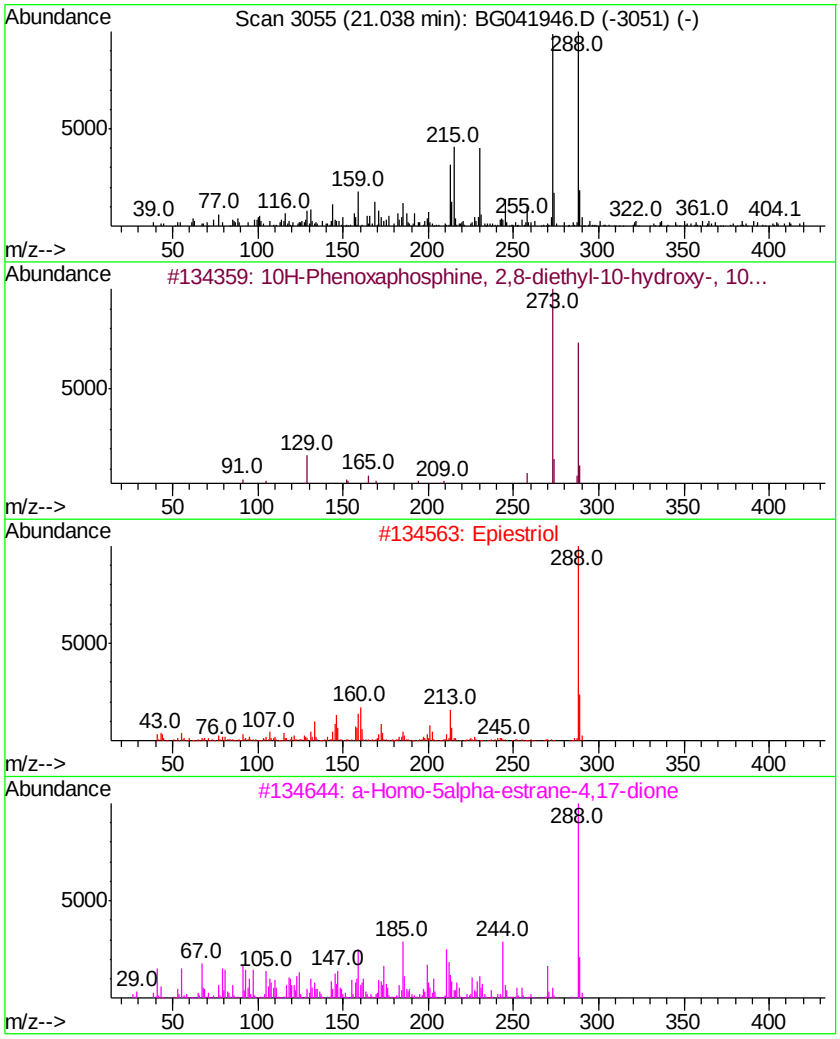
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

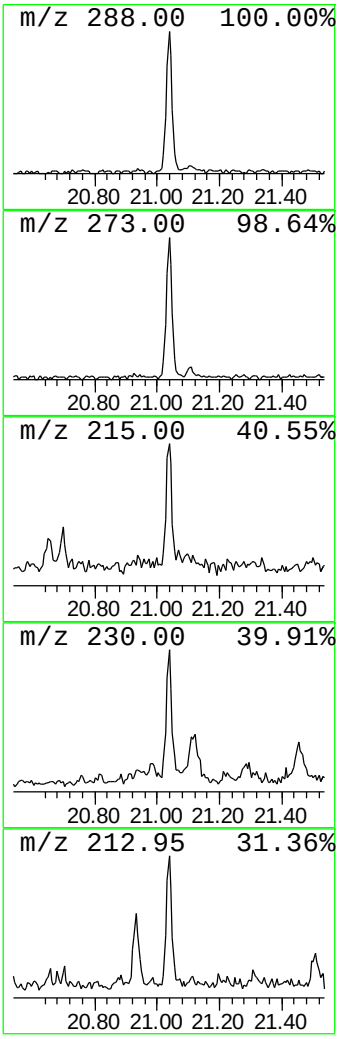
****

Peak Number 4 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	2.14 ng/ul	93940	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10H-Phenoxaphosphine, 2,8-diethy...	288	C16H17O3P	036360-92-6	47
2	Epiestriol	288	C18H24O3	000547-81-9	30
3	a-Homo-5alpha-estrane-4,17-dione	288	C19H28O2	014412-78-3	27
4	Pyrido[3,4-b]benzothiophene-4-ca...	273	C13H8FN3OS	1000265-50-3	25
5	1-Phosphacyclopentadiene, 2,3,4,...	288	C18H25OP	1000163-70-8	22





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

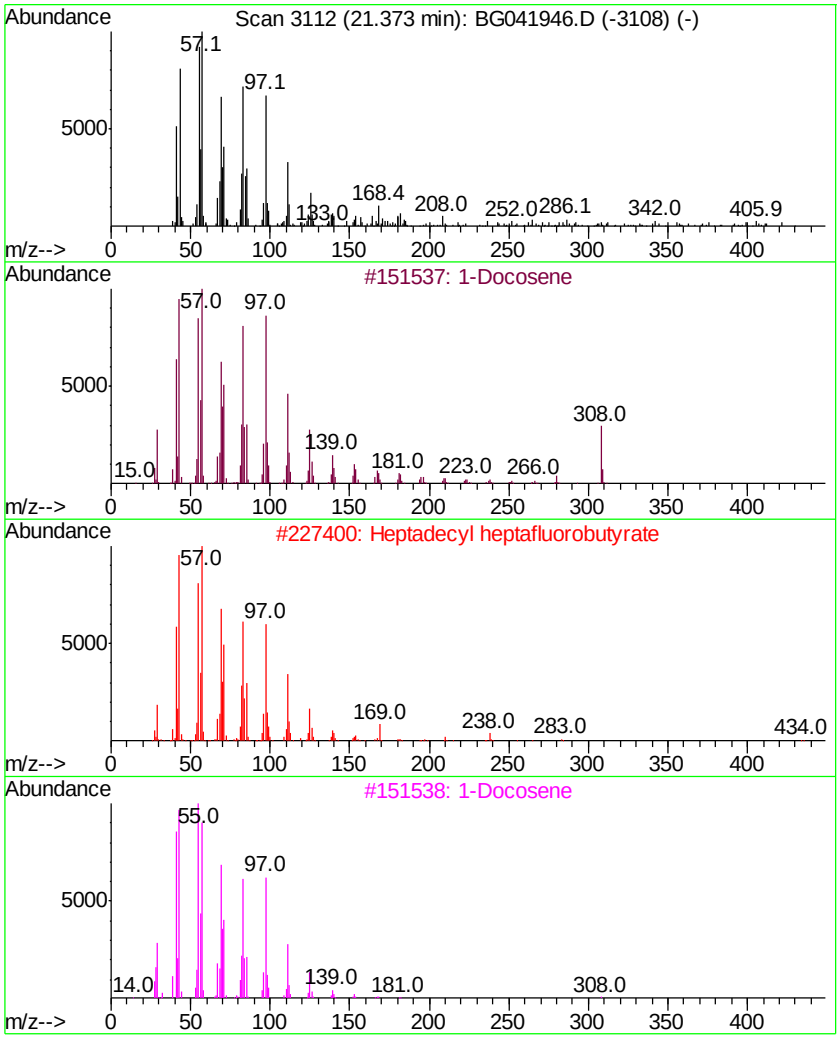
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

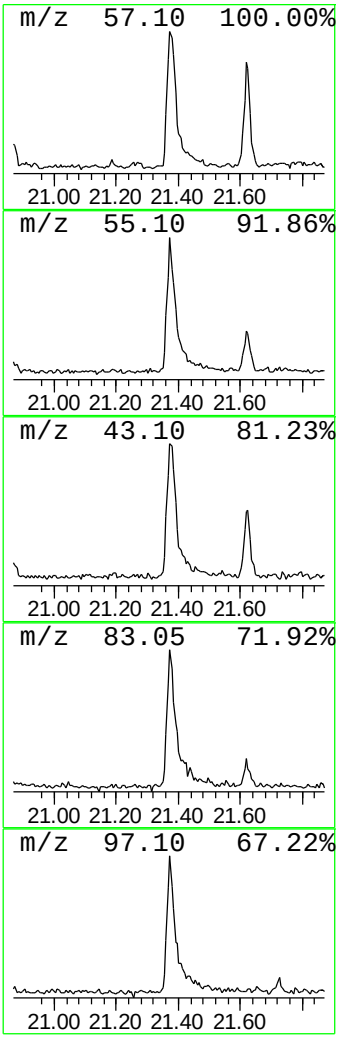
****

Peak Number 5 (DEL) Alkane: Cyclic21.37 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	4.29 ng/ul	188226	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	96
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3	1-Docosene	308	C22H44	001599-67-3	94
4	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	94
5	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	93





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Data File : BG041946.D
Acq On : 4 Aug 2019 17:36
Operator : HP/JU
Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

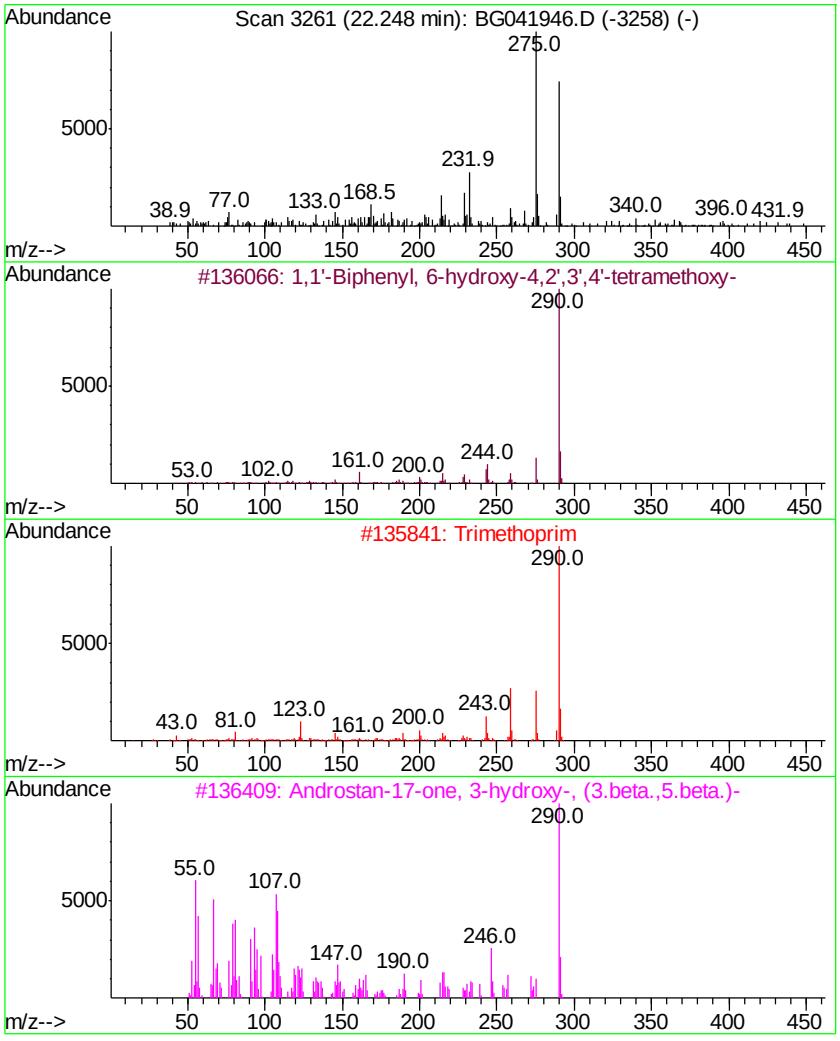
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

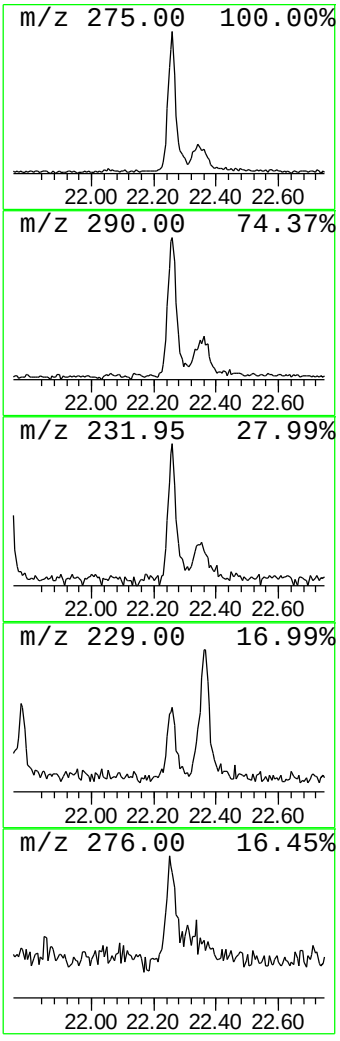
****

Peak Number 6 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.25	3.20 ng/ul	140165	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 6-hydroxy-4,2',3'...	290	C16H18O5	119101-16-5	46
2	Trimethoprim	290	C14H18N4O3	000738-70-5	43
3	Androstan-17-one, 3-hydroxy-, (3...	290	C19H30O2	000571-31-3	42
4	Benzo[ghi]perylene, 4-methyl-	290	C23H14	019224-38-5	38
5	Benzo[a]phenanthridin-4(1H)-one, ...	290	C19H18N2O	1000276-19-6	38





Library Search Compound Report

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Operator : HP/JU
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Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

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Quant Title : SVOA CALIBRATION

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

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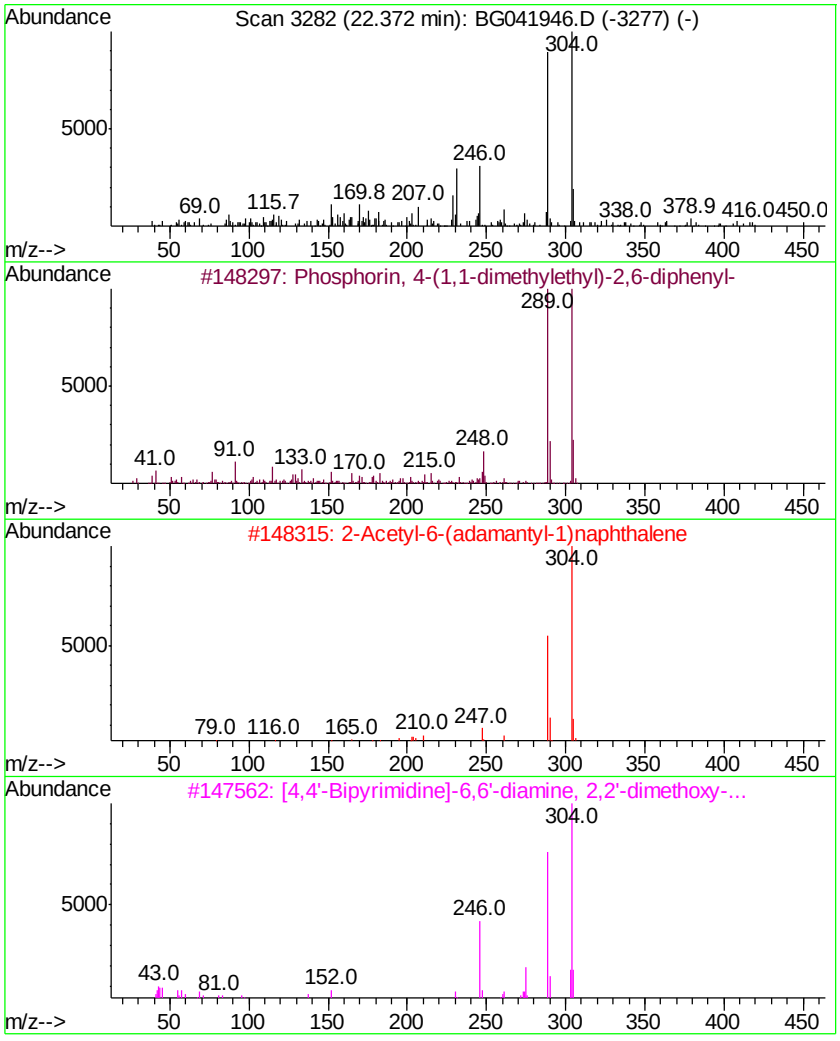
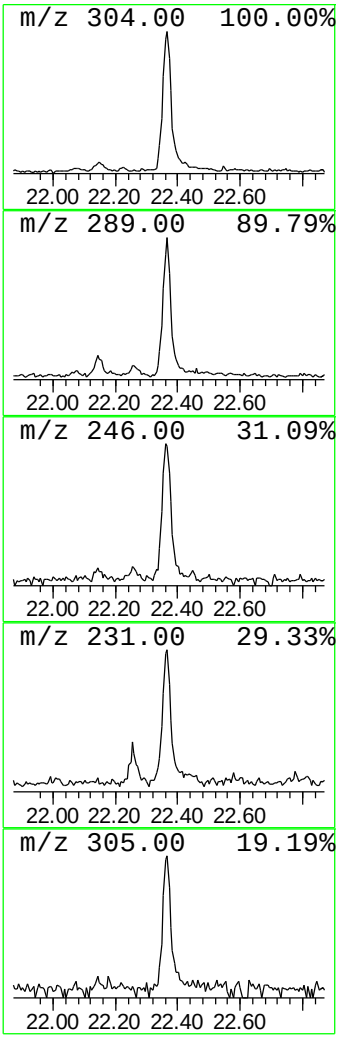
Peak Number 7 Phosphorin, 4-(1,1-dimethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	3.84 ng/ul	168334	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphorin, 4-(1,1-dimethylethyl...	304	C21H21P	017420-26-7	58
2	2-Acetyl-6-(adamantyl-1)naphthalene	304	C22H24O	066366-17-4	50
3	[4,4'-Bipyrimidine]-6,6'-diamine...	304	C14H20N6O2	062880-75-5	47
4	Imidazolo[2,1-b]thiazol-3-ol, 2,...	346	C11H11IN2OS	1000264-45-8	38
5	p-Diphosphorinane, 1,4-diphenyl-...	304	C16H18O2P2	109501-46-4	38





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Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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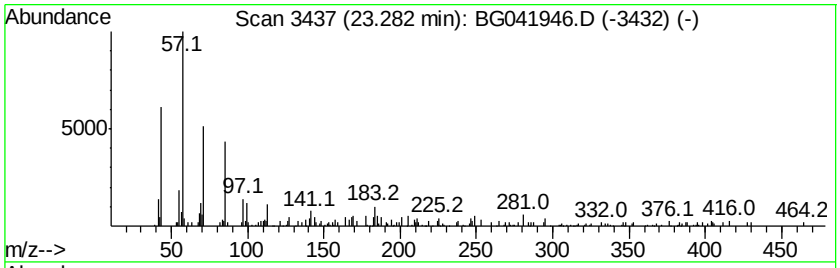
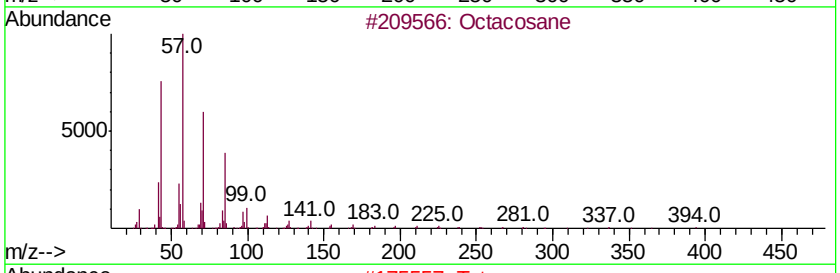
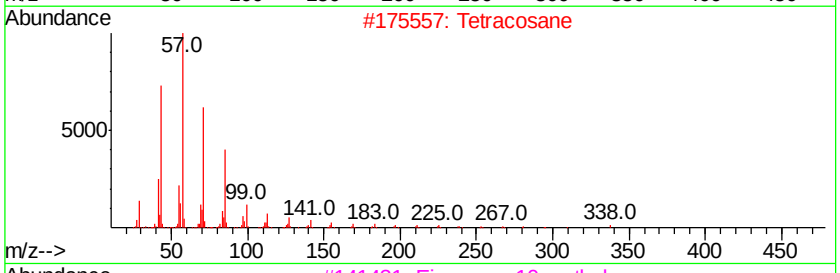
****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 8

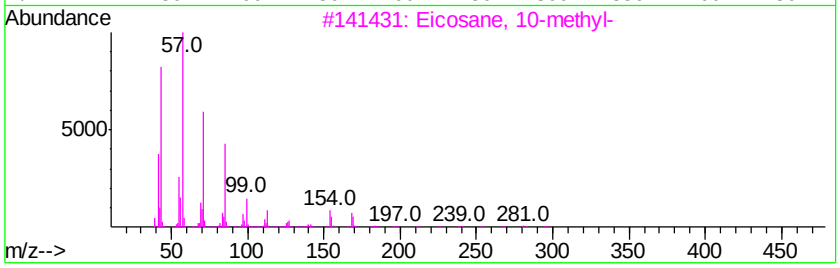
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.28	3.38 ng/ul	148158	Chrysene-d12	21.73

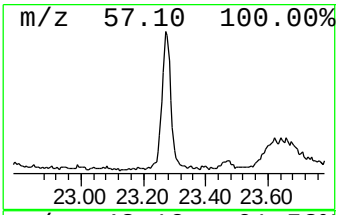
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	93
2	Tetracosane	338	C24H50	000646-31-1	91
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	64
4	Tricosane, 2-methyl-	338	C24H50	001928-30-9	58
5	Octacosane	394	C28H58	000630-02-4	58

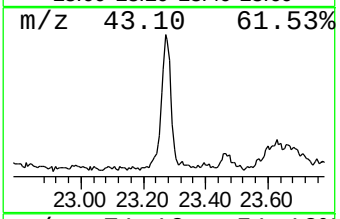


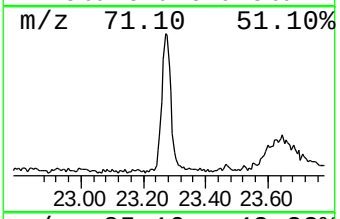


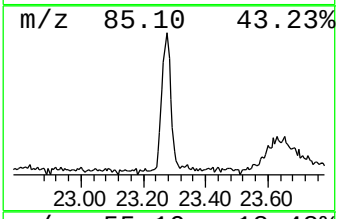


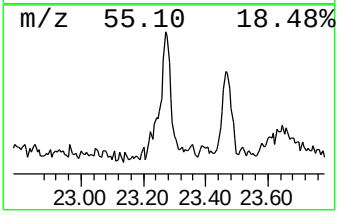












Library Search Compound Report

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Sample : K3962-15
Misc :
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Instrument : BNA_G
ClientSampleId : BEP08

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Quant Title : SVOA CALIBRATION

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

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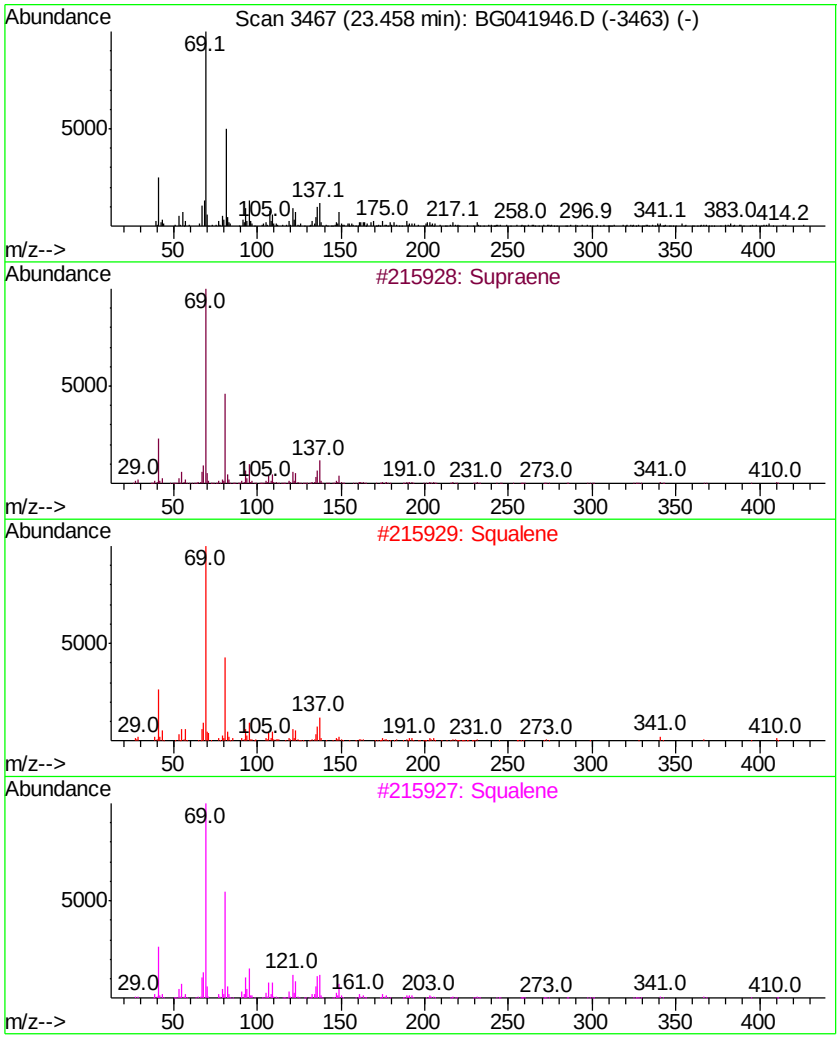
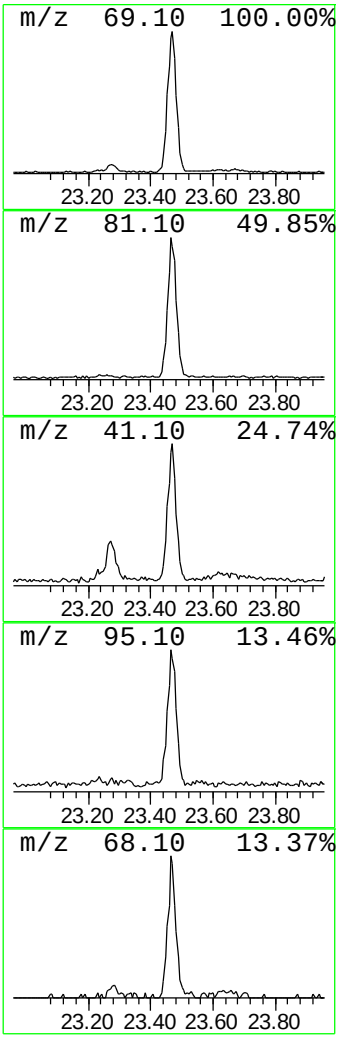
Peak Number 9 Supraene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.46	8.49 ng/ul	323536	Perylene-d12	25.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	96
2	Squalene	410	C30H50	000111-02-4	90
3	Squalene	410	C30H50	000111-02-4	87
4	Squalene	410	C30H50	000111-02-4	87
5	Squalene	410	C30H50	000111-02-4	87





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Quant Title : SVOA CALIBRATION

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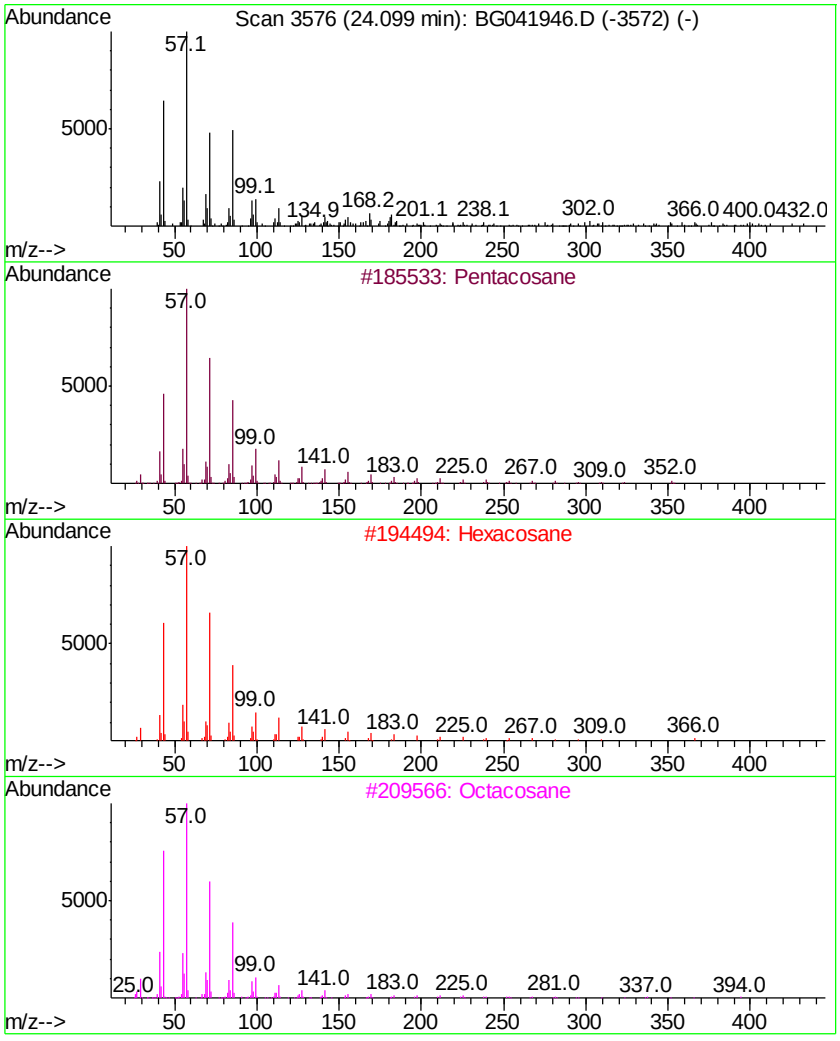
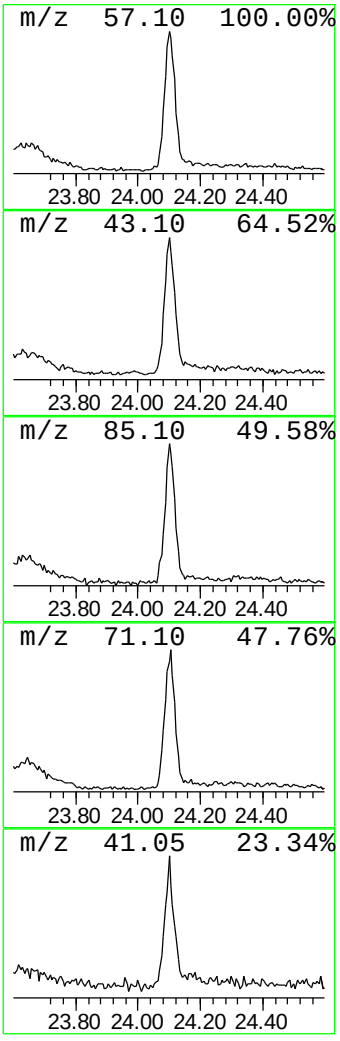
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.10	4.89 ng/ul	186157	Perylene-d12	25.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	93
2	Hexacosane	366	C26H54	000630-01-3	93
3	Octacosane	394	C28H58	000630-02-4	90
4	Hexacosane	366	C26H54	000630-01-3	87
5	Docosane	310	C22H46	000629-97-0	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080319\
Data File : BG041946.D
Acq On : 4 Aug 2019 17:36
Operator : HP/JU
Sample : K3962-15
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BEP08

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.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
Butane, 2-methoxy...	3.16	127.1	ng/ul	2044080	1	8.04	321581	20.0
n-Hexadecanoic acid	18.34	5.7	ng/ul	232726	4	17.45	813856	20.0
Androstane	20.50	3.6	ng/ul	158146	5	21.73	876808	20.0
unknown-01	21.04	2.1	ng/ul	93940	5	21.73	876808	20.0
(DEL) Alkane: Cyc...	21.37	4.3	ng/ul	188226	5	21.73	876808	20.0
unknown-02	22.25	3.2	ng/ul	140165	5	21.73	876808	20.0
Phosphorin, 4-(1,...	22.37	3.8	ng/ul	168334	5	21.73	876808	20.0
(DEL) Alkane: Str...	23.28	3.4	ng/ul	148158	5	21.73	876808	20.0
Supraene	23.46	8.5	ng/ul	323536	6	25.01	761892	20.0
(DEL) Alkane: Str...	24.10	4.9	ng/ul	186157	6	25.01	761892	20.0