LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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	3.146	5	10	32	rVB	1345248	2106786	90.02%	7.927%
2	3.410	51	55	62	rVB2	9769	15235	0.65%	0.057%
3	4.074	164	168	173	rVB3	9245	13027	0.56%	0.049%
4	7.170	689	695	706	rBV	159092	326711	13.96%	1.229%
5	7.341	715	724	735	rBV	137243	244319	10.44%	0.919%
6	7.552	753	760	775	rVV	188691	338694	14.47%	1.274%
7	8.028	833	841	851	rBV	185022	329828	14.09%	1.241%
8	8.733	953	961	973	rBV	193588	359692	15.37%	1.353%
9	9.192	1032	1039	1051	rBV	179783	333807	14.26%	1.256%
10	9.920	1155	1163	1173	rBV	157105	291719	12.46%	1.098%
11	10.467	1248	1256	1271	rBV	245713	482650	20.62%	1.816%
12	10.849	1313	1321	1329	rBV	281510	516159	22.05%	1.942%
13	10.978	1335	1343	1366	rBV	211805	434722	18.58%	1.636%
14	12.735	1636	1642	1650	rBV2	6178	11049	0.47%	0.042%
15	14.010	1853	1859	1864	rBV2	11396	17873	0.76%	0.067%
16	14.086	1864	1872	1876	rVV	590924	913929	39.05%	3.439%
17	14.133	1876	1880	1890	rVB	219751	335734	14.35%	1.263%
18	14.380	1915	1922	1933	rBV	685482	1094973	46.79%	4.120%
19	14.685	1966	1974	1981	rBV2	539020	847244	36.20%	3.188%
20	14.750	1981	1985	1995	rVB	22428	36144	1.54%	0.136%
21	14.867	1999	2005	2018	rBV	141363	275993	11.79%	1.038%
22	15.091	2038	2043	2051	rVB3	39661	63396	2.71%	0.239%
23	15.478	2104	2109	2113	rVV5	6487	13776	0.59%	0.052%
24	15.678	2134	2143	2149	rBV	948182	1455477	62.19%	5.476%
25	15.731	2149	2152	2156	rVB2	30692	42905	1.83%	0.161%
26	15.790	2156	2162	2168	rBV	196877	320263	13.68%	1.205%
27	15.919	2177	2184	2193	rVB4	11897	32849	1.40%	0.124%
28	16.031	2198	2203	2209	rVB2	13534	22180	0.95%	0.083%
29	16.178	2220	2228	2235	rVB	16207	33302	1.42%	0.125%
30	16.407	2263	2267	2274	rBV5	6610	12253	0.52%	0.046%
31	16.724	2315	2321	2322	rVV4	7744	12449	0.53%	0.047%
32	16.742	2322	2324	2329	rVV4	11289	15010	0.64%	0.056%
33	16.789	2329	2332	2337	rVB3	6734	10911	0.47%	0.041%
34	16.977	2355	2364	2367	rBV5	11289	29034	1.24%	0.109%

LSC Area Percent Report

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Sample : K4029-16
Misc :
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Instrument : BNA_G
ClientSampleId : BENY6

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

35	17.012	2367	2370	2378	rVV4	11840	19116	0.82%	0.072%
36	17.088	2378	2383	2390	rVV3	10875	20135	0.86%	0.076%
37	17.171	2392	2397	2405	rVV7	11196	27082	1.16%	0.102%
38	17.259	2409	2412	2421	rVB4	9272	16301	0.70%	0.061%
39	17.435	2435	2442	2447	rBV	721380	1172102	50.08%	4.410%
40	17.476	2447	2449	2454	rVV	206702	285782	12.21%	1.075%
41	17.535	2454	2459	2470	rVV	1043886	1713175	73.20%	6.446%
42	17.623	2470	2474	2480	rVB6	11016	20582	0.88%	0.077%
43	17.840	2501	2511	2515	rBV2	16169	38158	1.63%	0.144%
44	18.028	2538	2543	2550	rVB7	6032	13304	0.57%	0.050%
45	18.322	2586	2593	2596	rVV2	50230	103453	4.42%	0.389%
46	18.363	2596	2600	2606	rVV	73571	134728	5.76%	0.507%
47	18.446	2610	2614	2619	rVV	34444	54489	2.33%	0.205%
48	18.510	2619	2625	2629	rVV3	54442	111890	4.78%	0.421%
49	18.545	2629	2631	2637	rVB	30980	41576	1.78%	0.156%
50	18.622	2639	2644	2648	rBV8	6592	10921	0.47%	0.041%
51	18.786	2668	2672	2674	rVV	30751	43277	1.85%	0.163%
52	18.810	2674	2676	2680	rVV	24792	37247	1.59%	0.140%
53	18.851	2680	2683	2688	rVV2	13580	19959	0.85%	0.075%
54	18.939	2695	2698	2704	rVB7	6477	11169	0.48%	0.042%
55	19.057	2714	2718	2723	rBV5	21593	28338	1.21%	0.107%
56	19.145	2730	2733	2739	rVB4	9765	13401	0.57%	0.050%
57	19.227	2739	2747	2750	rBV2	31067	59571	2.55%	0.224%
58	19.292	2755	2758	2762	rBV3	12116	14514	0.62%	0.055%
59	19.368	2767	2771	2778	rVB6	18446	38280	1.64%	0.144%
60	19.491	2780	2792	2797	rBV	225301	363659	15.54%	1.368%
61	19.544	2799	2801	2805	rVB5	10732	12904	0.55%	0.049%
62	19.585	2805	2808	2813	rBV5	12480	18502	0.79%	0.070%
63	19.685	2821	2825	2826	rBV3	8401	10878	0.46%	0.041%
64	19.826	2842	2849	2861	rBV2	1372957	2340342	100.00%	8.806%
65	19.926	2861	2866	2872	rVB4	25745	49680	2.12%	0.187%
66	20.026	2878	2883	2889	rBV2	20631	42373	1.81%	0.159%
67	20.085	2890	2893	2900	rVB4	14258	24297	1.04%	0.091%
68	20.167	2903	2907	2915	rBV2	28239	81256	3.47%	0.306%
69	20.308	2926	2931	2932	rVV5	19369	26389	1.13%	0.099%
70	20.343	2933	2937	2943	rVB2	51927	86432	3.69%	0.325%
71	20.443	2950	2954	2957	rBV2	29670	36014	1.54%	0.136%

LSC Area Percent Report

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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

72	20.496	2957	2963	2968	rVB3	38240	68293	2.92%	0.257%
73	20.584	2975	2978	2983	rVB6	10602	16066	0.69%	0.060%
74	20.643	2983	2988	2991	rBV	22026	32861	1.40%	0.124%
75	20.690	2991	2996	2999	rVV2	23525	40246	1.72%	0.151%
76	20.719	2999	3001	3006	rVB6	11240	16669	0.71%	0.063%
77	20.790	3011	3013	3018	rVB4	10517	15430	0.66%	0.058%
78	20.854	3022	3024	3027	rVV3	13739	13534	0.58%	0.051%
79	20.937	3028	3038	3048	rVV3	42765	183556	7.84%	0.691%
80	21.107	3063	3067	3073	rVB2	23559	45986	1.96%	0.173%
81	21.366	3103	3111	3119	rBV5	89255	228793	9.78%	0.861%
82	21.436	3121	3123	3134	rVB	34123	72127	3.08%	0.271%
83	21.718	3162	3171	3176	rBV2	863421	1516516	64.80%	5.706%
84	21.765	3176	3179	3187	rVB	90041	146536	6.26%	0.551%
85	21.924	3202	3206	3212	rVB3	37777	57368	2.45%	0.216%
86	22.129	3237	3241	3248	rVB5	13099	28454	1.22%	0.107%
87	22.547	3306	3312	3318	rBV3	44420	88738	3.79%	0.334%
88	22.729	3341	3343	3349	rVB4	14932	19610	0.84%	0.074%
89	23.258	3428	3433	3440	rVB2	57111	107964	4.61%	0.406%
90	23.951	3542	3551	3558	rBV3	50832	183207	7.83%	0.689%
91	24.080	3566	3573	3581	rBV	69576	160910	6.88%	0.605%
92	24.674	3668	3674	3678	rBV2	22356	55023	2.35%	0.207%
93	24.756	3678	3688	3710	rVB	719567	2152485	91.97%	8.099%
94	24.985	3716	3727	3734	rBV	492668	1364391	58.30%	5.134%
95	25.055	3736	3739	3749	rVB2	73125	156393	6.68%	0.588%
96	26.213	3929	3936	3947	rVB3	61287	193804	8.28%	0.729%
97	27.253	4102	4113	4144	rBV9	139800	770404	32.92%	2.899%
98	27.605	4166	4173	4186	rVB3	53031	171763	7.34%	0.646%
99	29.286	4453	4459	4471	rVB10	28817	110576	4.72%	0.416%
100	31.307	4796	4803	4804	rBV6	15328	27844	1.19%	0.105%

Sum of corrected areas: 26576916

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
Data File : BG042019.D
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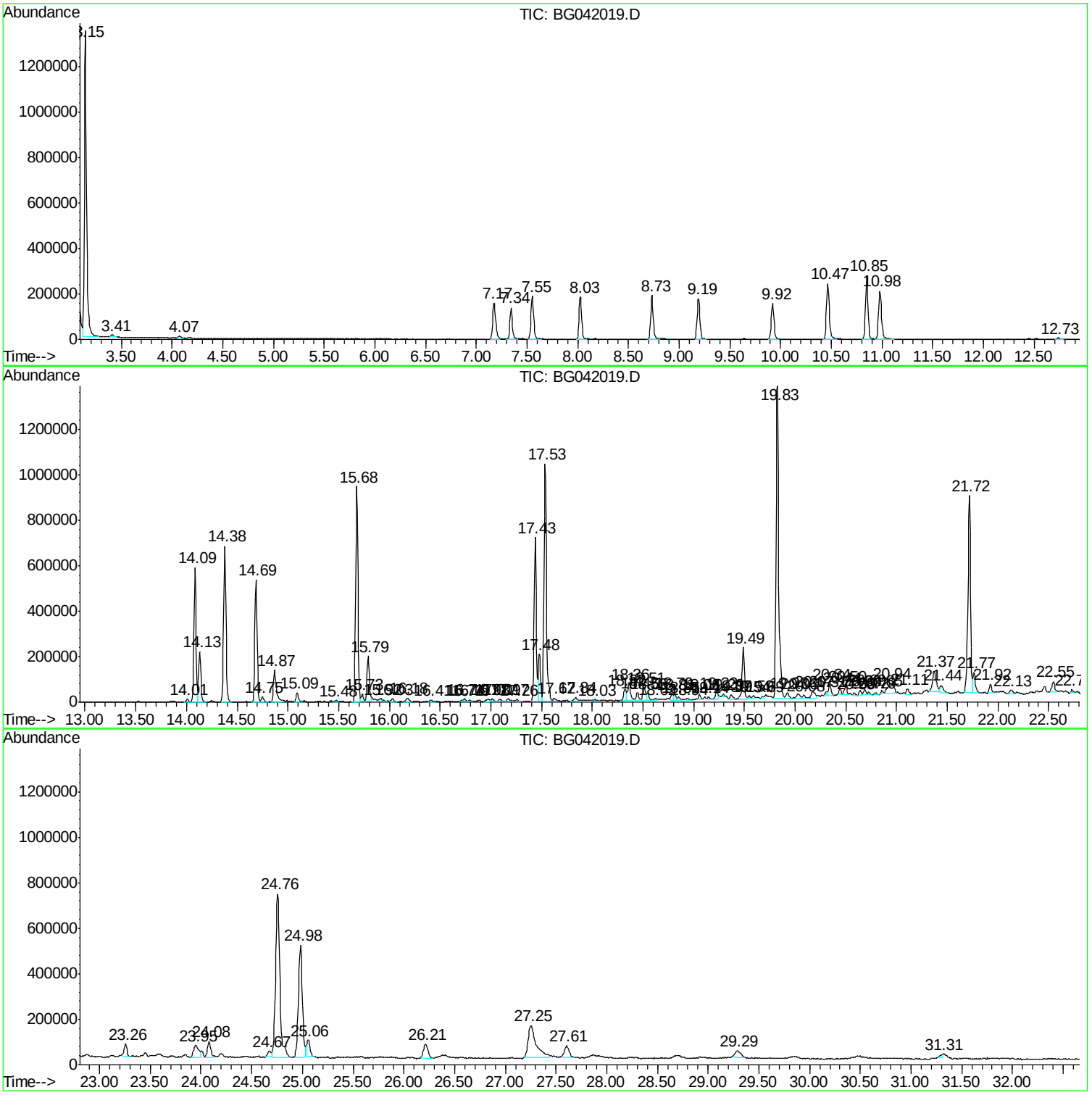
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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\BG080619\
Data File : BG042019.D
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Instrument : BNA_G
ClientSampleId : BENY6

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

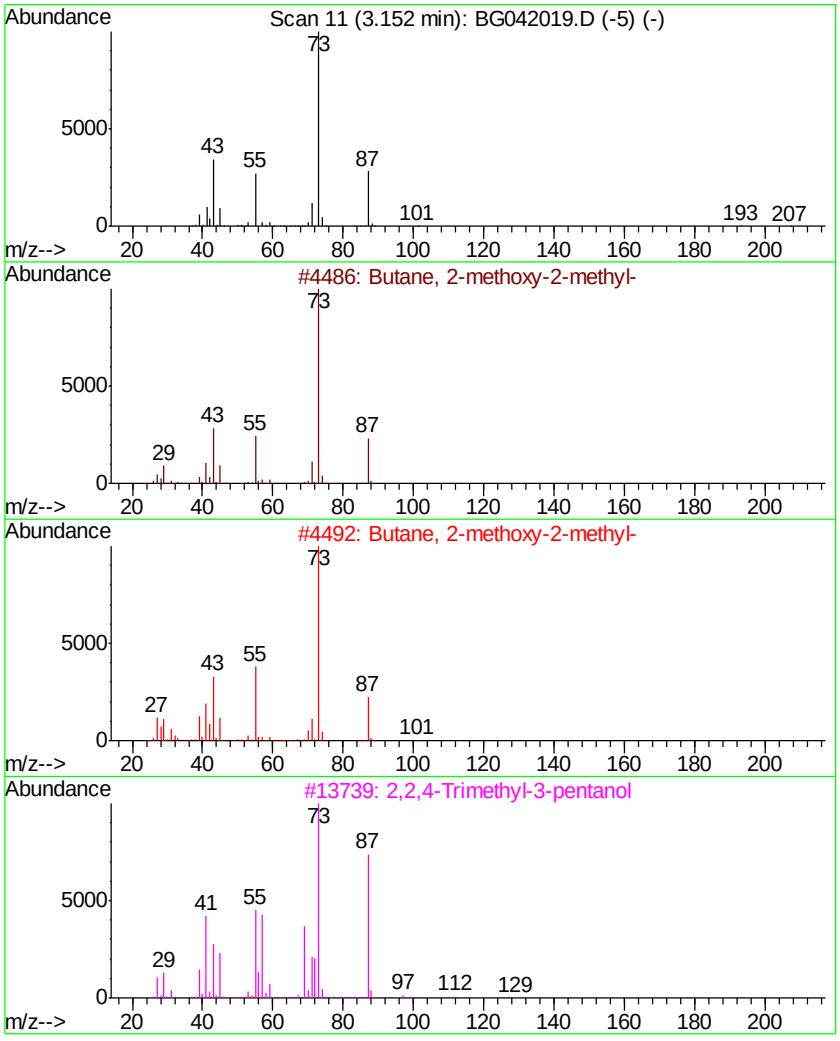
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

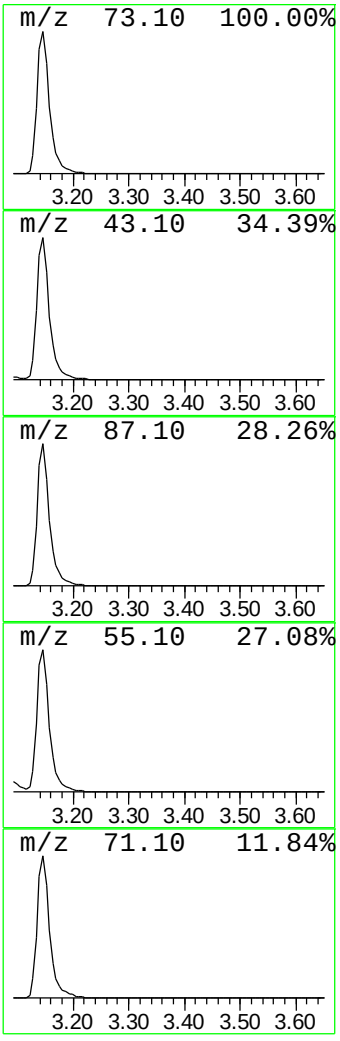
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	127.75 ng/ul	2106790	1,4-Dichlorobenzene-d4	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
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Misc :
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Instrument : BNA_G
ClientSampleId : BENY6

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

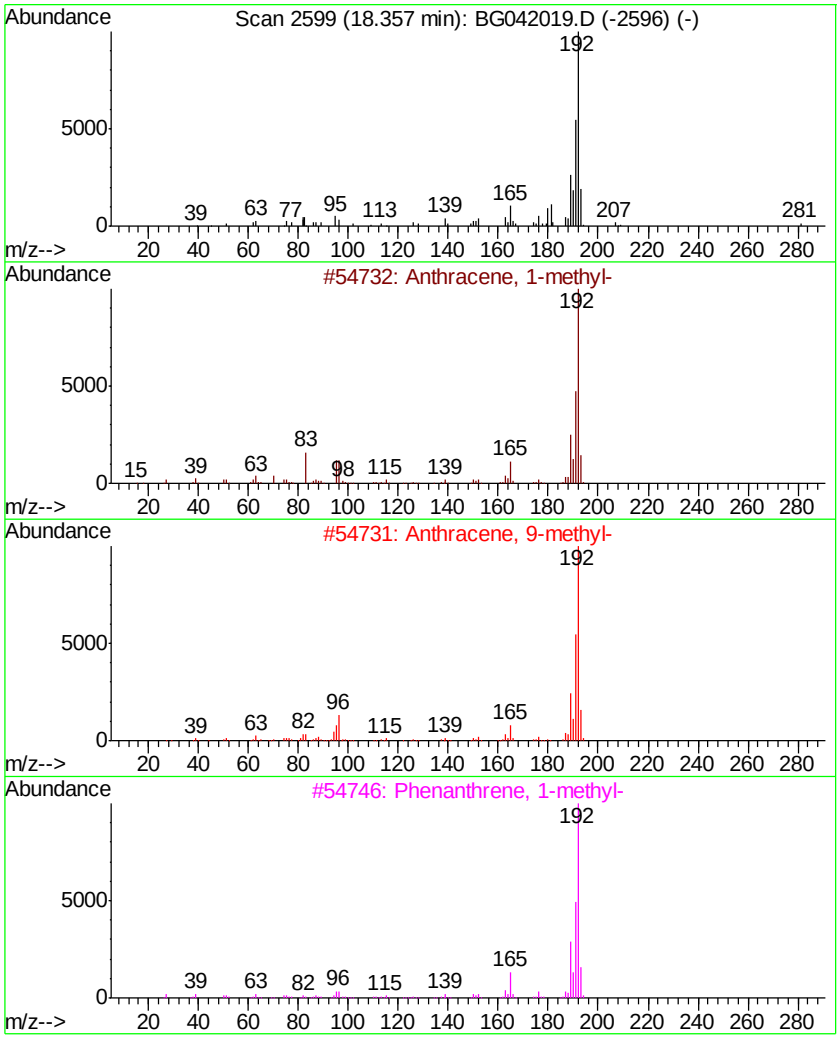
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

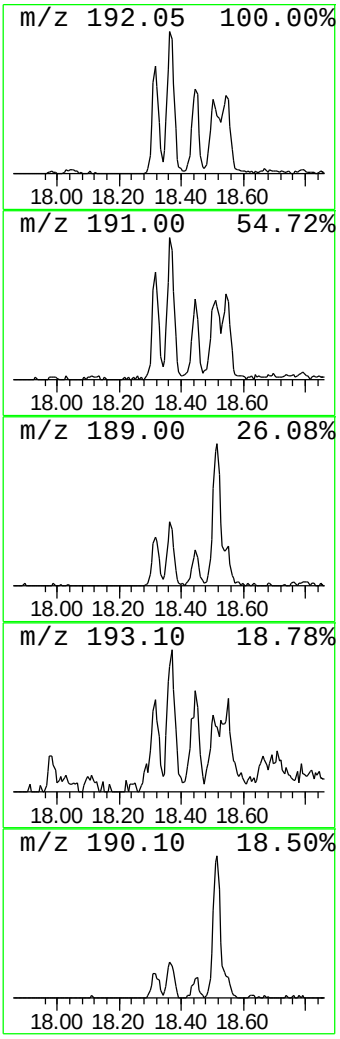
****

Peak Number 2 Anthracene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.30 ng/ul	134728	Phenanthrene-d10	17.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87





Library Search Compound Report

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Instrument : BNA_G
ClientSampleId : BENY6

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

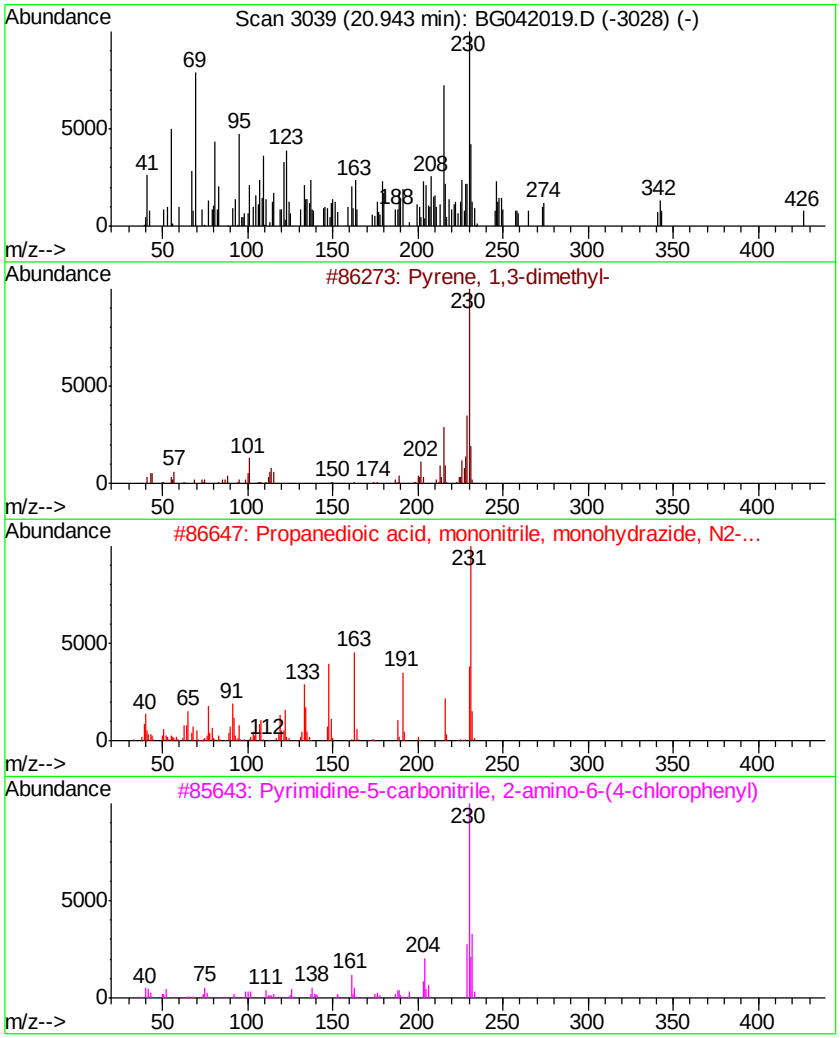
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

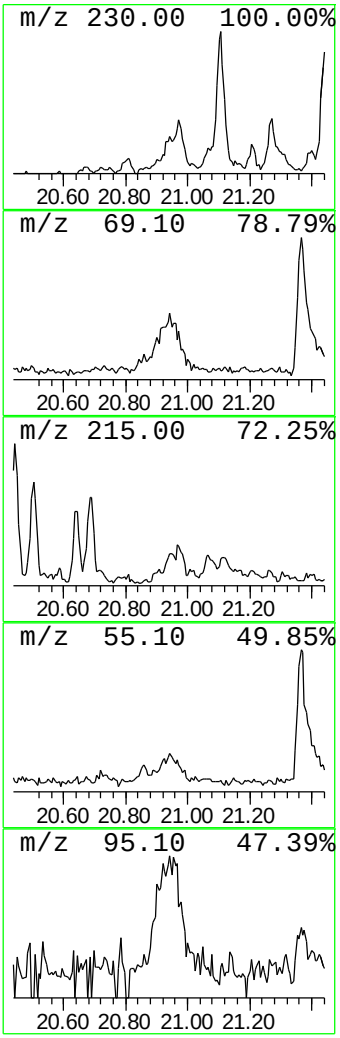
****

Peak Number 3 Pyrene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	2.42 ng/ul	183556	Chrysene-d12	21.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	60
2	Propanedioic acid, mononitrile, ...	231	C12H13N3O2	004980-37-4	35
3	Pyrimidine-5-carbonitrile, 2-ami...	230	C11H7ClN4	281665-63-2	35
4	Pyrene, 1-(methoxymethyl)-	246	C18H14O	091385-15-8	35
5	2'-Acetonaphthone, 1',8'-dihydro...	230	C14H14O3	000838-03-9	30





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Data File : BG042019.D
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Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

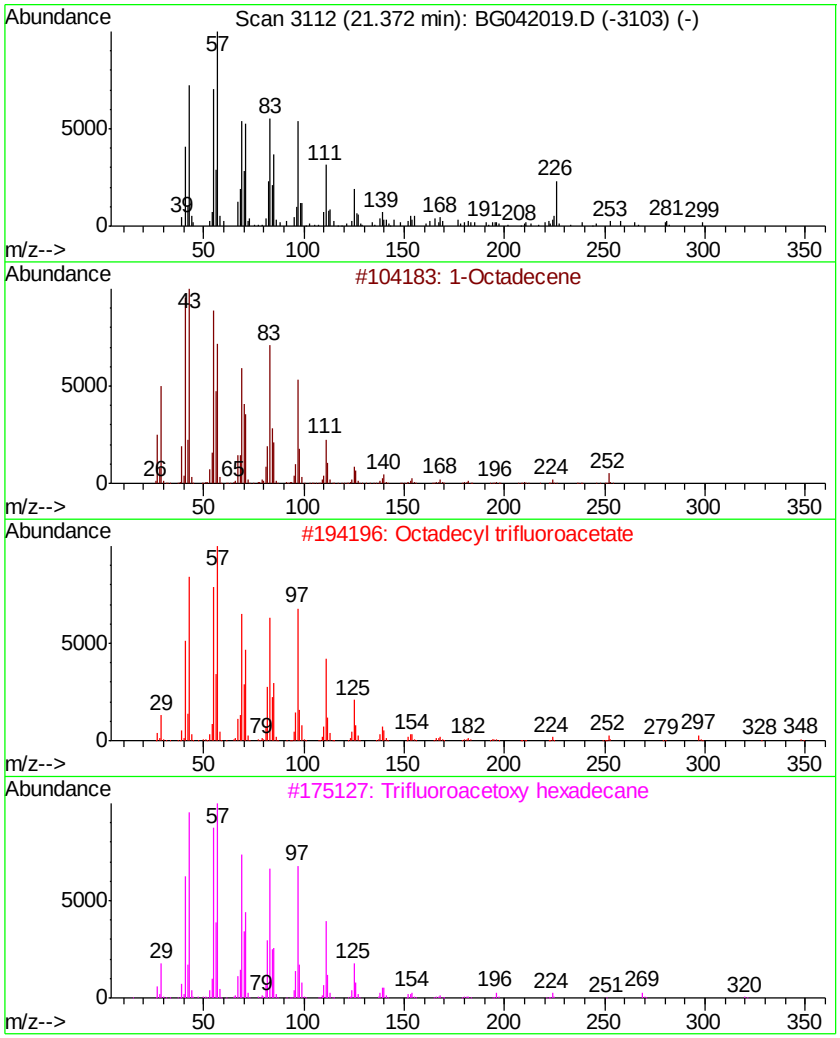
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

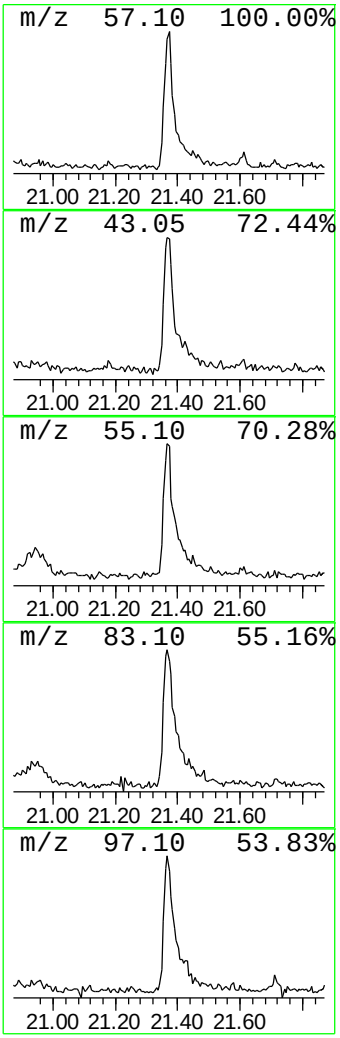
****

Peak Number 4 (DEL) Alkane: Cyclic21.37 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	3.02 ng/ul	228793	Chrysene-d12	21.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	97
2	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	93
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
4	1-Octadecene	252	C18H36	000112-88-9	93
5	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91





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Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

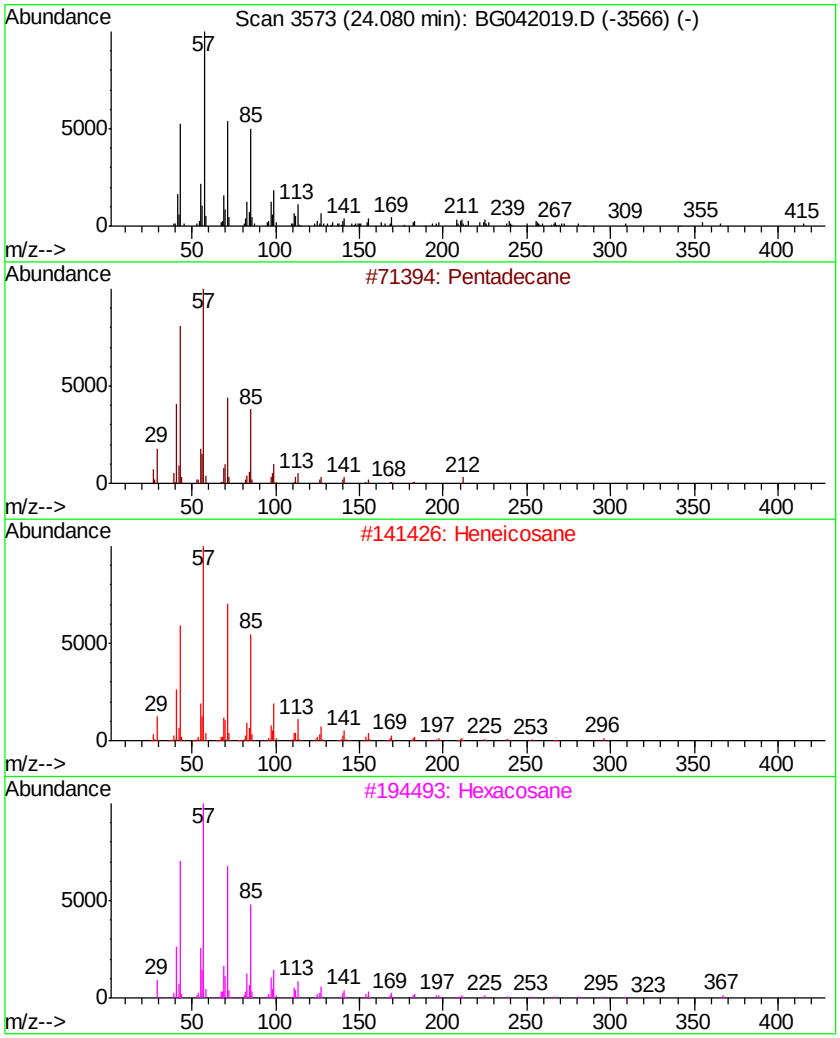
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

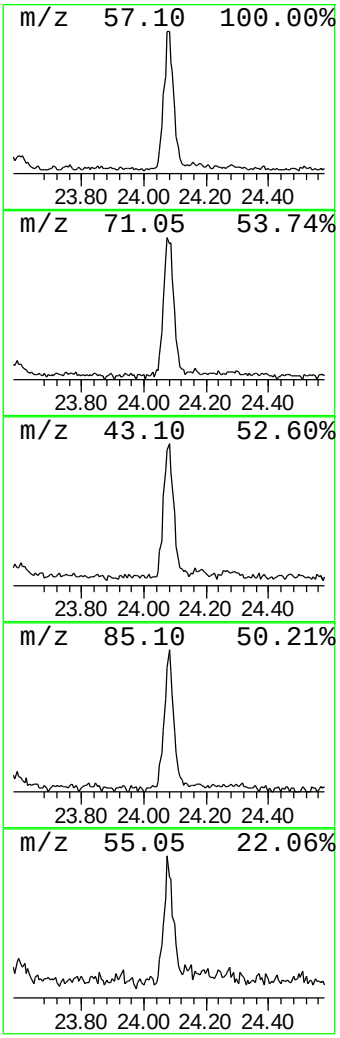
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	2.36 ng/ul	160910	Perylene-d12	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Hexacosane	366	C26H54	000630-01-3	90
4	Tetracosane	338	C24H50	000646-31-1	90
5	Hentriacontane	437	C31H64	000630-04-6	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

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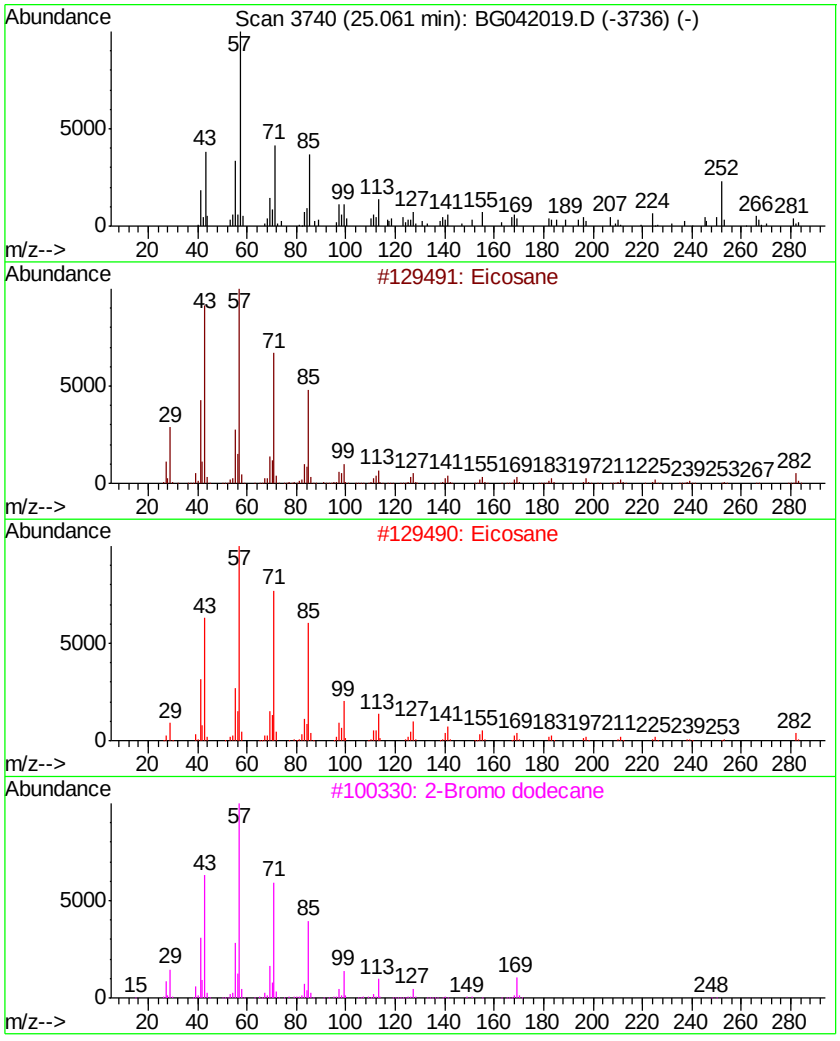
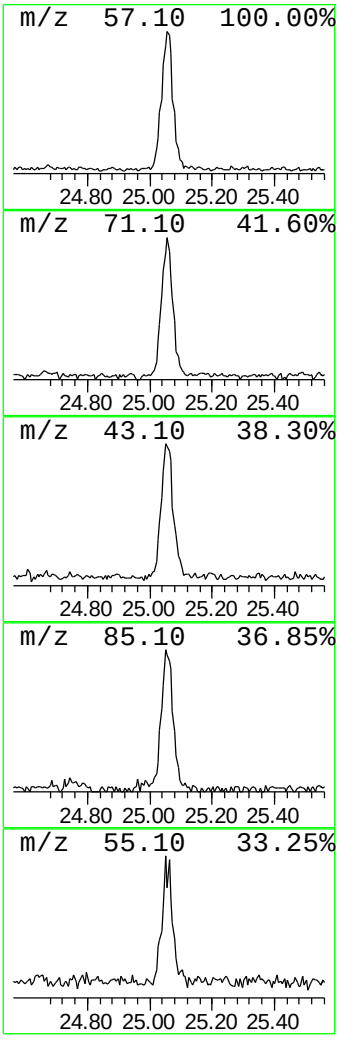
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.06	2.29 ng/ul	156393	Perylene-d12	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	78
2	Eicosane	282	C20H42	000112-95-8	78
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	49
4	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	49
5	Heptadecane	240	C17H36	000629-78-7	47





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

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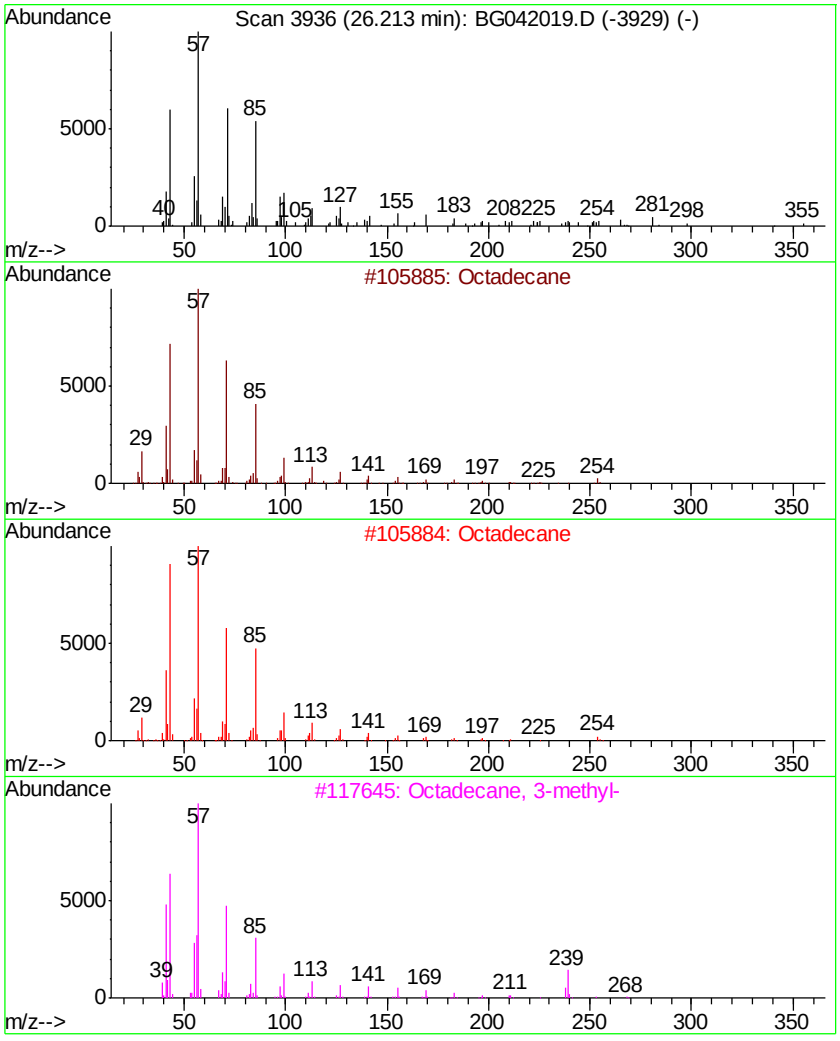
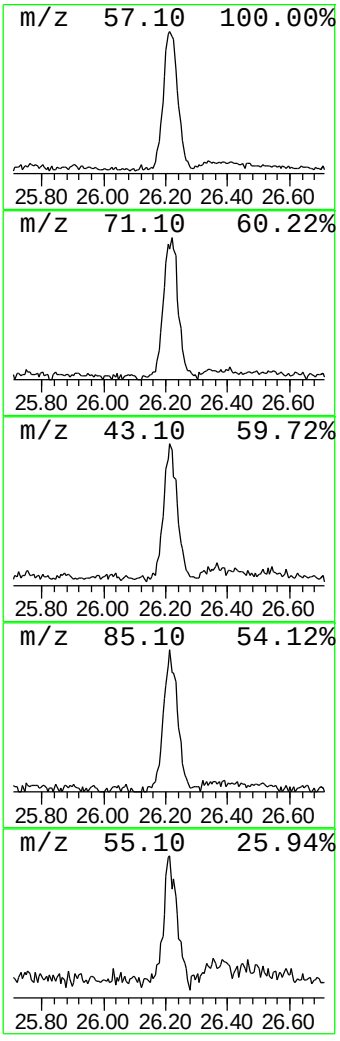
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.21	2.84 ng/ul	193804	Perylene-d12	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	94
2	Octadecane	254	C18H38	000593-45-3	91
3	Octadecane, 3-methyl-	268	C19H40	006561-44-0	89
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	87
5	Octacosane	394	C28H58	000630-02-4	87





Library Search Compound Report

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Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

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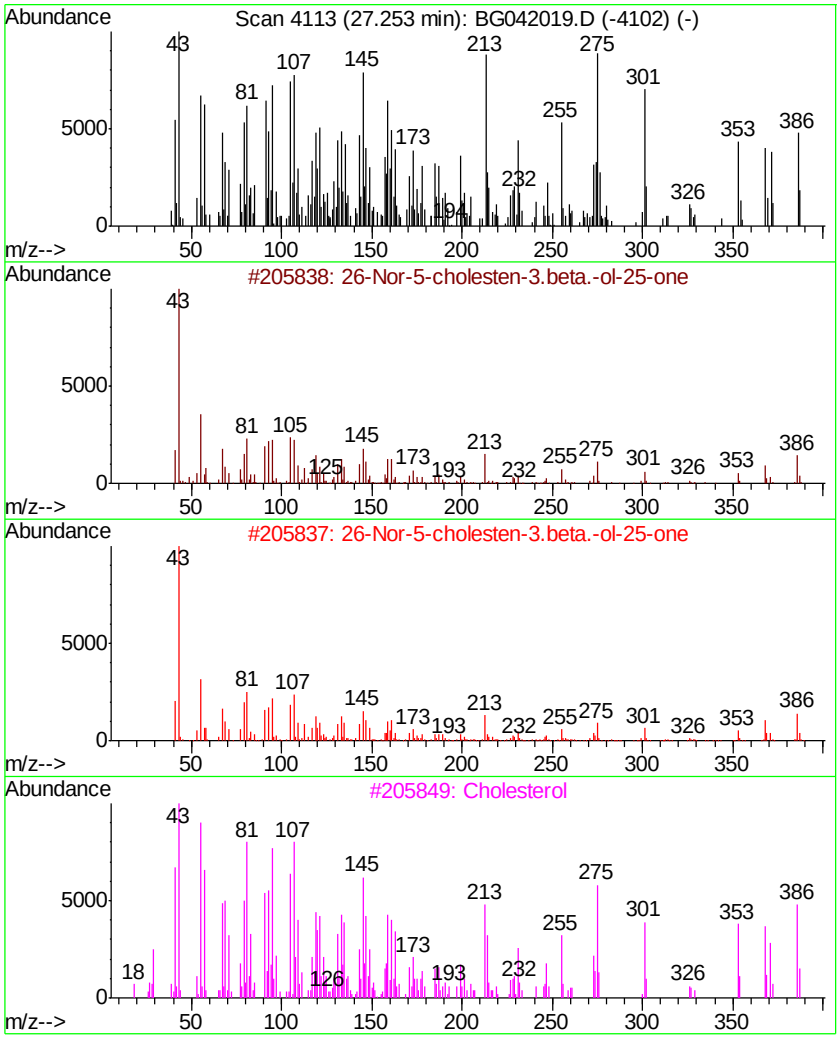
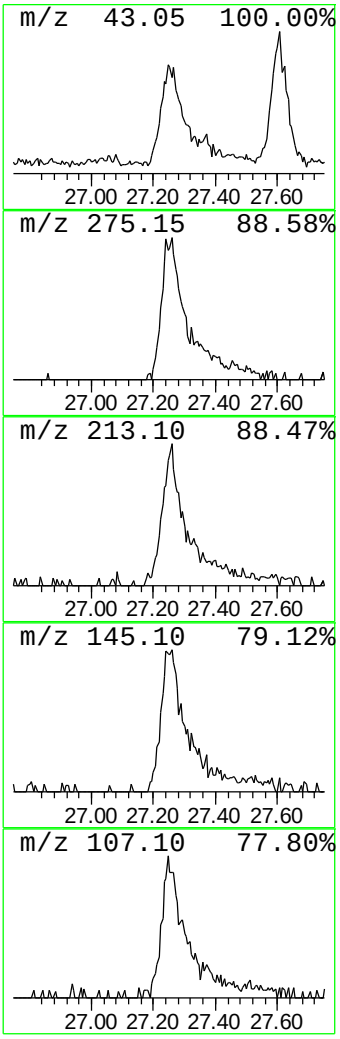
Peak Number 8 26-Nor-5-cholesten-3.beta.-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.25	11.29 ng/ul	770404	Perylene-d12	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
2	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
3	Cholesterol	386	C27H46O	000057-88-5	96
4	Cholesterol	386	C27H46O	000057-88-5	93
5	17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-50-0	15





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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

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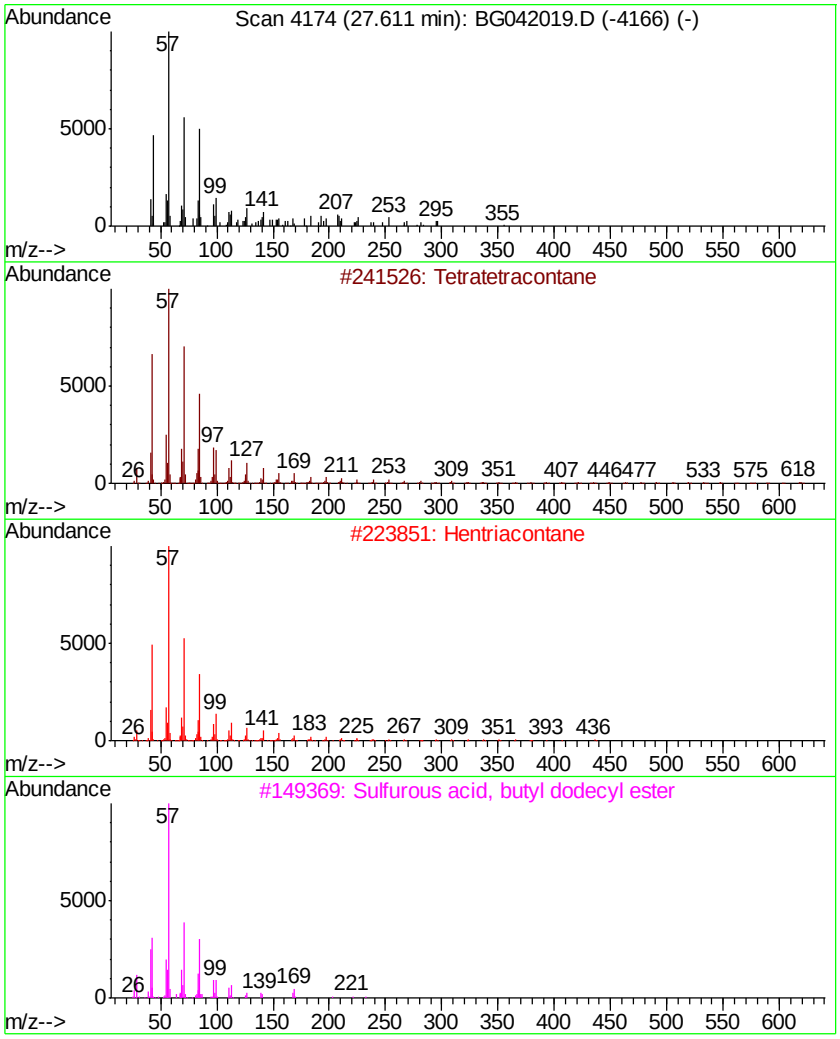
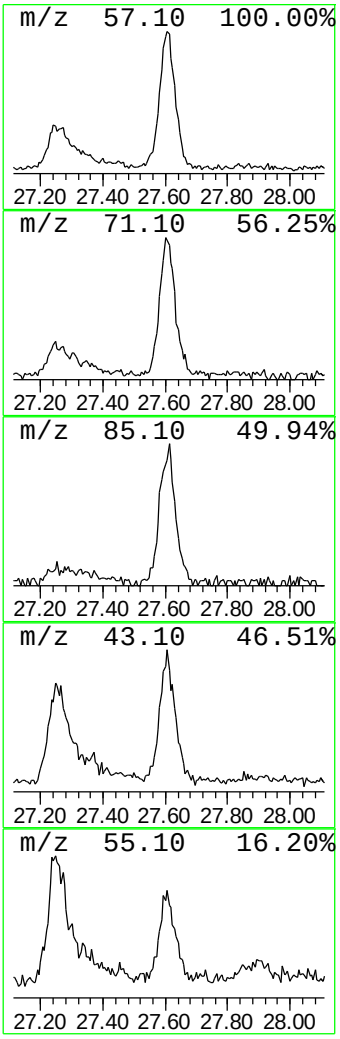
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.61	2.52 ng/ul	171763	Perylene-d12	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	91
2	Hentriacontane	437	C31H64	000630-04-6	87
3	Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	86
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	86
5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080619\
Data File : BG042019.D
Acq On : 7 Aug 2019 5:20
Operator : HP/JU
Sample : K4029-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BENY6

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...	3.15	127.8	ng/ul	2106790	1	8.03	329828	20.0
Anthracene, 1-met...	18.36	2.3	ng/ul	134728	4	17.43	1172100	20.0
Pyrene, 1,3-dimet...	20.94	2.4	ng/ul	183556	5	21.72	1516520	20.0
(DEL) Alkane: Cyc...	21.37	3.0	ng/ul	228793	5	21.72	1516520	20.0
(DEL) Alkane: Str...	24.08	2.4	ng/ul	160910	6	24.98	1364390	20.0
(DEL) Alkane: Str...	25.06	2.3	ng/ul	156393	6	24.98	1364390	20.0
(DEL) Alkane: Str...	26.21	2.8	ng/ul	193804	6	24.98	1364390	20.0
26-Nor-5-choleste...	27.25	11.3	ng/ul	770404	6	24.98	1364390	20.0
(DEL) Alkane: Str...	27.61	2.5	ng/ul	171763	6	24.98	1364390	20.0