

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
 Data File : BG019281.D
 Acq On : 21 Oct 2015 14:52
 Operator : UM/NP
 Sample : G4074-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9K1

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG101515.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	3.078	36	41	48	rVV	16206	27336	5.83%	0.426%
2	3.154	49	54	64	rVB	32003	45381	9.69%	0.707%
3	3.419	94	99	105	rVB2	1920	3056	0.65%	0.048%
4	5.093	380	384	389	rBV3	1429	1895	0.40%	0.030%
5	6.997	706	708	715	rVB2	725	958	0.20%	0.015%
6	7.161	729	736	744	rBV	59764	107447	22.93%	1.673%
7	7.308	755	761	768	rBV	40591	66367	14.16%	1.033%
8	7.520	789	797	805	rBV	60300	100995	21.55%	1.572%
9	7.984	870	876	885	rVB	72740	118452	25.28%	1.844%
10	8.701	991	998	1007	rVV	77284	134671	28.74%	2.097%
11	9.147	1067	1074	1082	rVB	65174	107640	22.97%	1.676%
12	9.277	1093	1096	1104	rBV4	1996	3882	0.83%	0.060%
13	9.870	1191	1197	1204	rBV	57365	101095	21.58%	1.574%
14	10.416	1283	1290	1303	rBV	87859	152534	32.55%	2.375%
15	10.792	1347	1354	1363	rBV	116866	198422	42.35%	3.089%
16	10.928	1371	1377	1384	rBV	61821	112381	23.98%	1.750%
17	11.586	1486	1489	1495	rBV2	1389	2225	0.47%	0.035%
18	12.373	1619	1623	1627	rVB2	1121	1563	0.33%	0.024%
19	12.978	1720	1726	1731	rVB3	2263	3081	0.66%	0.048%
20	13.225	1763	1768	1772	rVB4	3702	4921	1.05%	0.077%
21	14.024	1897	1904	1908	rBV	177281	252825	53.96%	3.936%
22	14.071	1908	1912	1918	rVB	76870	105013	22.41%	1.635%
23	14.265	1941	1945	1946	rVV2	882	981	0.21%	0.015%
24	14.312	1947	1953	1961	rVV	181389	284891	60.80%	4.436%
25	14.506	1982	1986	1990	rBV	1997	2436	0.52%	0.038%
26	14.623	1998	2006	2012	rBV2	196279	314718	67.17%	4.900%
27	14.841	2035	2043	2053	rBV	81180	133162	28.42%	2.073%
28	15.011	2069	2072	2073	rBV3	803	943	0.20%	0.015%
29	15.029	2073	2075	2080	rVB3	1516	2024	0.43%	0.032%
30	15.416	2133	2141	2144	rVV2	2732	4880	1.04%	0.076%
31	15.458	2144	2148	2152	rVB2	7102	9854	2.10%	0.153%
32	15.499	2152	2155	2157	rBV	1120	1140	0.24%	0.018%
33	15.610	2168	2174	2180	rBV	252973	376984	80.45%	5.869%
34	15.657	2180	2182	2186	rVB3	849	992	0.21%	0.015%

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35	15.734	2190	2195	2203	rVB	88332	126472	26.99%	1.969%
36	15.851	2211	2215	2221	rVB5	3436	6866	1.47%	0.107%
37	16.010	2237	2242	2245	rBV3	1530	1819	0.39%	0.028%
38	16.257	2281	2284	2285	rVB	1065	979	0.21%	0.015%
39	16.280	2285	2288	2289	rBV	905	1096	0.23%	0.017%
40	16.321	2291	2295	2297	rBV2	12299	15800	3.37%	0.246%
41	16.398	2307	2308	2311	rBV2	1192	933	0.20%	0.015%
42	16.498	2322	2325	2329	rVB2	4949	6042	1.29%	0.094%
43	16.592	2339	2341	2344	rVB2	894	1108	0.24%	0.017%
44	16.627	2344	2347	2350	rBV3	637	1051	0.22%	0.016%
45	16.656	2350	2352	2357	rBV4	1098	1831	0.39%	0.029%
46	16.756	2367	2369	2373	rBV3	2150	2789	0.60%	0.043%
47	16.827	2378	2381	2382	rBV3	1466	1145	0.24%	0.018%
48	16.897	2389	2393	2395	rBV4	1958	2984	0.64%	0.046%
49	16.926	2397	2398	2401	rVV3	2153	1722	0.37%	0.027%
50	16.991	2407	2409	2412	rBV3	1382	1579	0.34%	0.025%
51	17.038	2415	2417	2420	rVB3	1378	1359	0.29%	0.021%
52	17.109	2425	2429	2435	rVV2	11224	16056	3.43%	0.250%
53	17.167	2435	2439	2443	rVB5	4882	7498	1.60%	0.117%
54	17.261	2451	2455	2457	rBV4	2520	3141	0.67%	0.049%
55	17.320	2461	2465	2467	rVB4	911	1313	0.28%	0.020%
56	17.367	2467	2473	2479	rBV	260434	386681	82.52%	6.020%
57	17.420	2479	2482	2483	rVV3	1920	1949	0.42%	0.030%
58	17.467	2483	2490	2498	rVB	265997	409653	87.43%	6.378%
59	17.532	2498	2501	2502	rBV2	2591	1696	0.36%	0.026%
60	17.584	2507	2510	2515	rVB5	1194	1751	0.37%	0.027%
61	17.637	2515	2519	2520	rBV3	1638	1496	0.32%	0.023%
62	17.737	2533	2536	2540	rVV3	2559	4991	1.07%	0.078%
63	17.773	2540	2542	2545	rVV4	2410	2657	0.57%	0.041%
64	17.849	2551	2555	2559	rVV2	8628	12189	2.60%	0.190%
65	17.878	2559	2560	2562	rVV	1690	1285	0.27%	0.020%
66	17.902	2562	2564	2566	rVV3	1697	1628	0.35%	0.025%
67	17.949	2570	2572	2573	rVV2	1458	1154	0.25%	0.018%
68	17.978	2576	2577	2580	rVV2	1884	1708	0.36%	0.027%
69	18.013	2581	2583	2584	rVV	1638	1436	0.31%	0.022%
70	18.031	2584	2586	2588	rVV3	2478	1957	0.42%	0.030%
71	18.060	2588	2591	2592	rVB3	1154	923	0.20%	0.014%

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72	18.125	2600	2602	2603	rBV	3932	2853	0.61%	0.044%
73	18.196	2609	2614	2619	rBV2	27699	39675	8.47%	0.618%
74	18.254	2619	2624	2634	rBV2	101334	144615	30.86%	2.252%
75	18.354	2639	2641	2644	rVB4	2781	2716	0.58%	0.042%
76	18.466	2656	2660	2661	rVB4	2187	1998	0.43%	0.031%
77	18.542	2670	2673	2675	rBV3	5943	5868	1.25%	0.091%
78	18.589	2679	2681	2682	rBV2	2030	1094	0.23%	0.017%
79	18.671	2692	2695	2697	rBV4	3497	4561	0.97%	0.071%
80	18.765	2709	2711	2714	rBV4	1757	1976	0.42%	0.031%
81	18.807	2714	2718	2721	rVV6	2393	4362	0.93%	0.068%
82	18.871	2728	2729	2733	rBV4	2775	3198	0.68%	0.050%
83	18.918	2735	2737	2740	rVV4	2732	2533	0.54%	0.039%
84	18.995	2748	2750	2755	rBV6	3050	4638	0.99%	0.072%
85	19.036	2755	2757	2758	rBV	2921	1766	0.38%	0.027%
86	19.106	2766	2769	2770	rBV2	4229	3955	0.84%	0.062%
87	19.177	2778	2781	2785	rVB6	3519	4529	0.97%	0.071%
88	19.206	2785	2786	2788	rBV2	2405	2100	0.45%	0.033%
89	19.253	2791	2794	2796	rBV4	3113	3885	0.83%	0.060%
90	19.382	2813	2816	2818	rBV4	8638	10559	2.25%	0.164%
91	19.523	2836	2840	2848	rBV3	31073	50849	10.85%	0.792%
92	19.753	2873	2879	2886	rBV	319886	468567	100.00%	7.295%
93	20.246	2961	2963	2964	rBV2	6888	4966	1.06%	0.077%
94	21.274	3133	3138	3150	rVB	92096	146801	31.33%	2.286%
95	21.633	3193	3199	3205	rVB	278903	450828	96.21%	7.019%
96	21.809	3227	3229	3230	rBV2	12159	9689	2.07%	0.151%
97	22.420	3329	3333	3346	rVB	88498	170376	36.36%	2.653%
98	23.906	3581	3586	3593	rVB	63491	122149	26.07%	1.902%
99	24.606	3697	3705	3714	rVB	169921	464553	99.14%	7.233%
100	24.829	3735	3743	3754	rVB	165818	465347	99.31%	7.245%

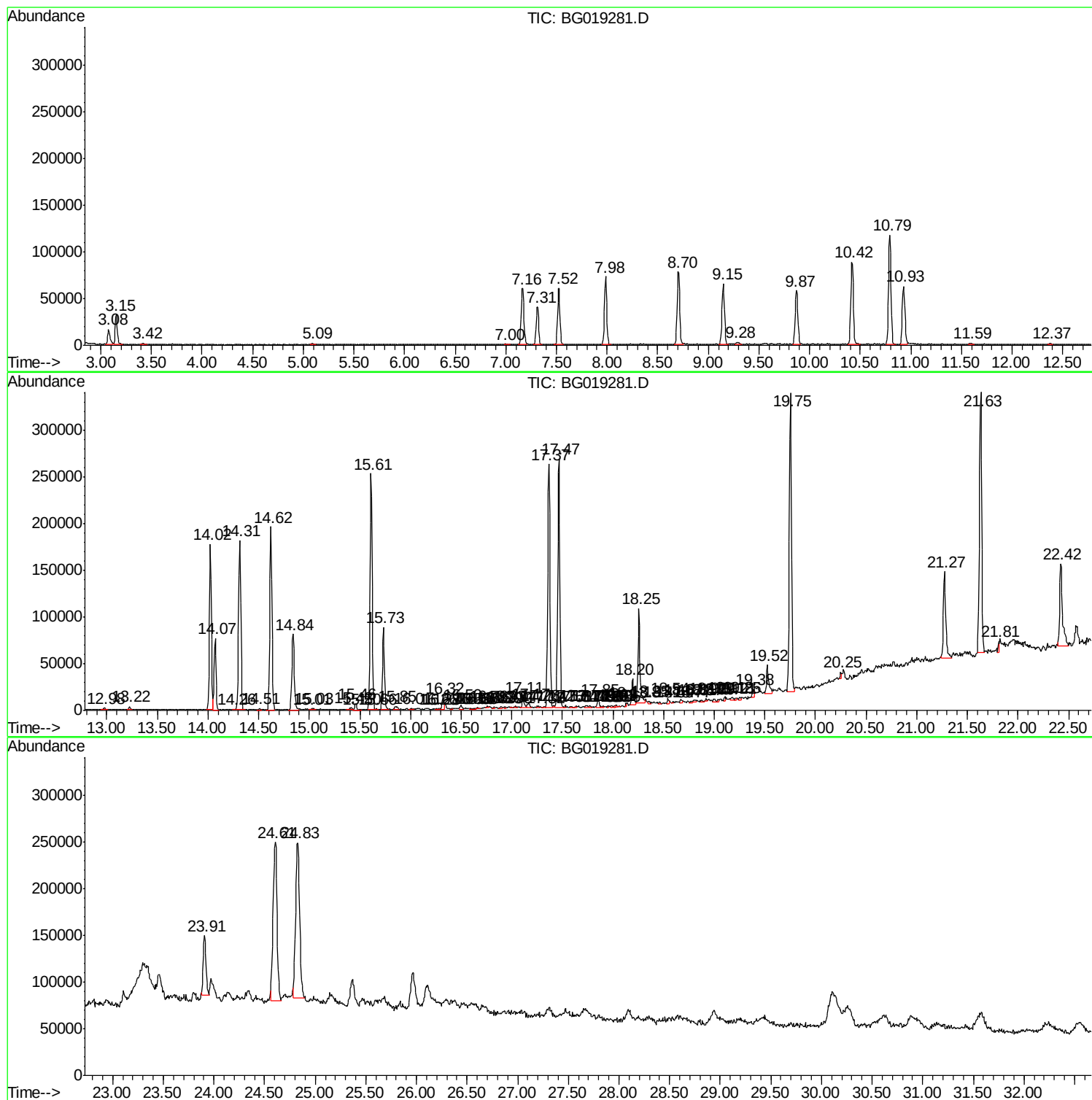
Sum of corrected areas: 6422888

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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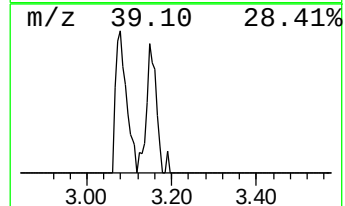
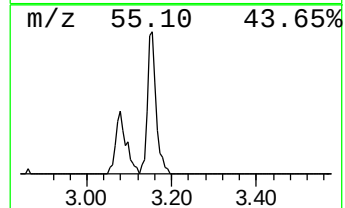
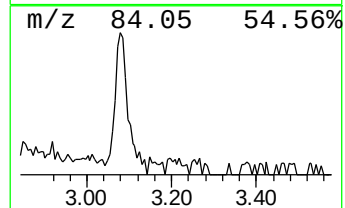
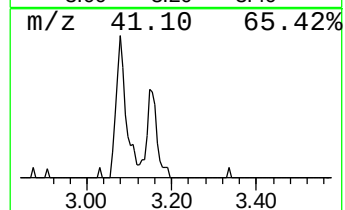
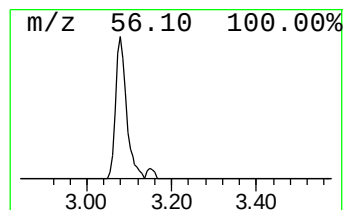
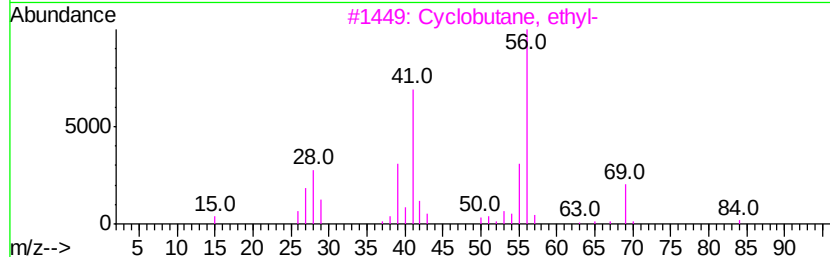
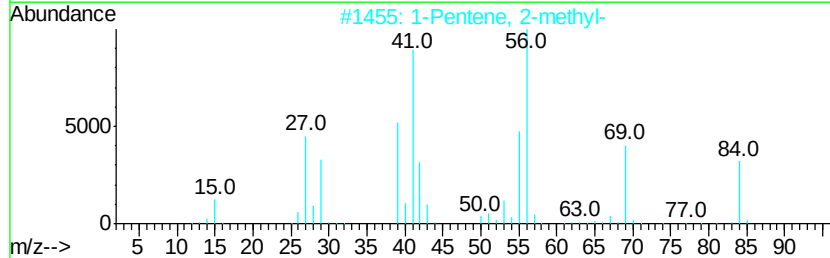
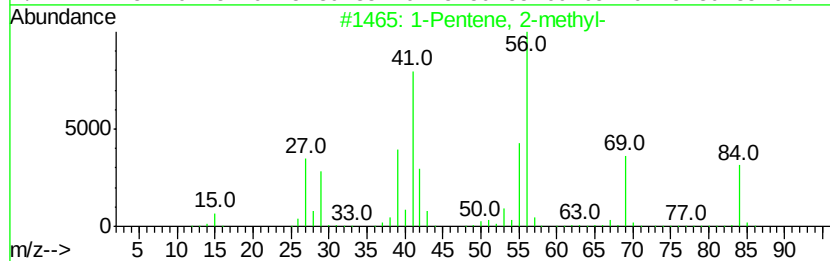
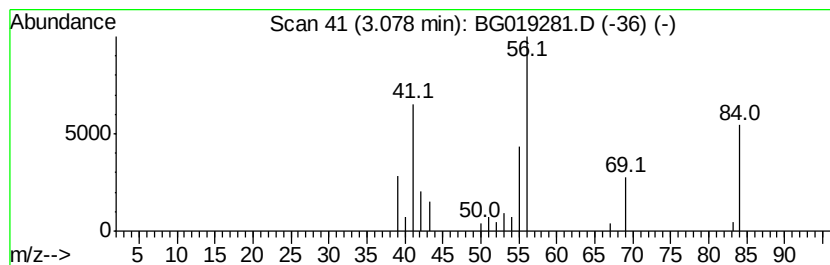
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.08	4.62 ng/ul	27336	1,4-Dichlorobenzene-d4	7.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
2			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	50
4			1-Hexene, 3,4-dimethyl-	112	C8H16	016745-94-1	39
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	38



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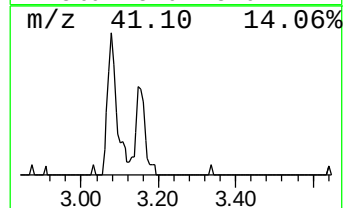
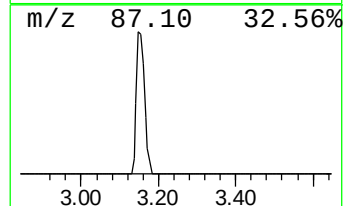
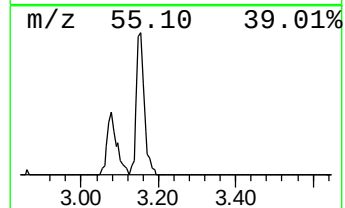
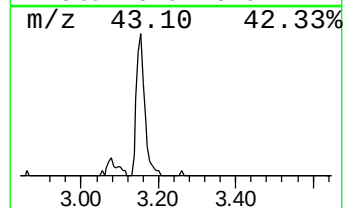
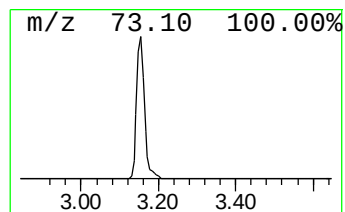
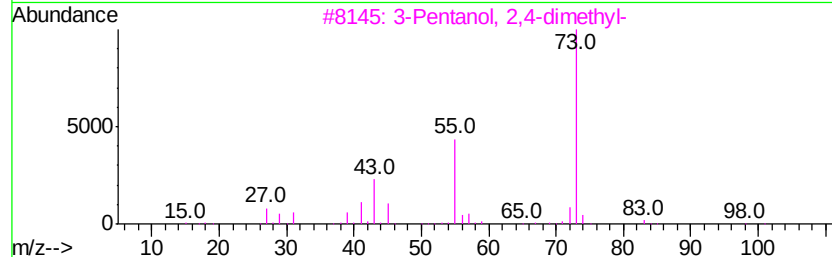
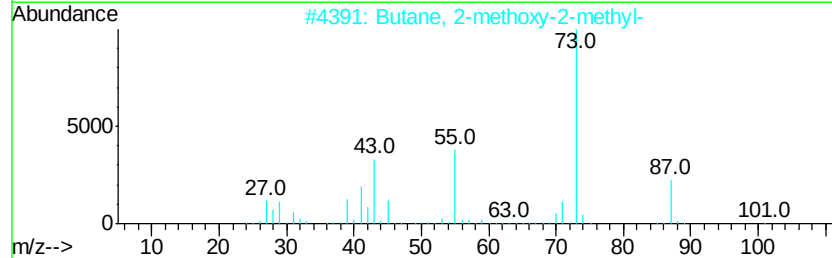
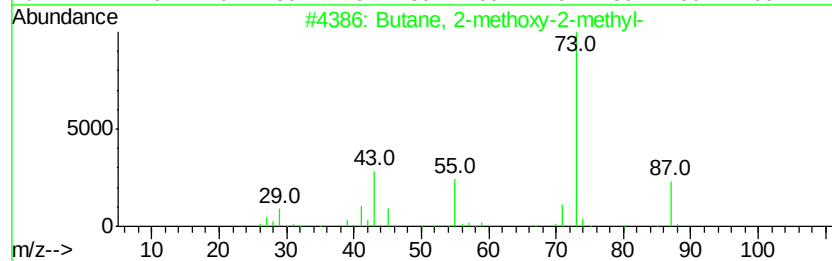
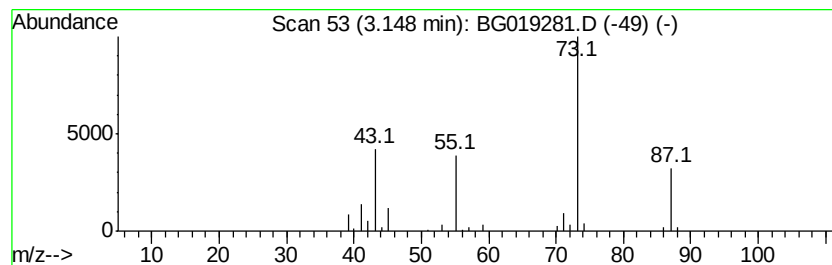
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	7.66 ng/ul	45381	1,4-Dichlorobenzene-d4	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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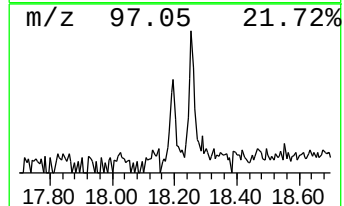
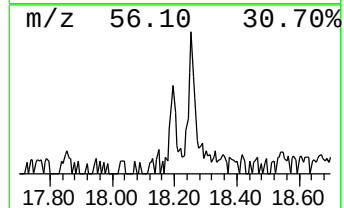
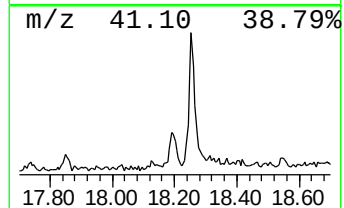
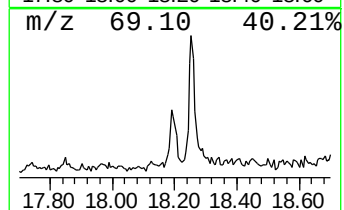
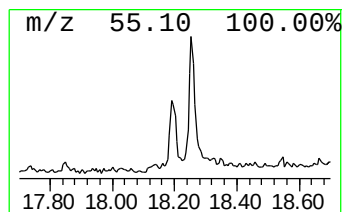
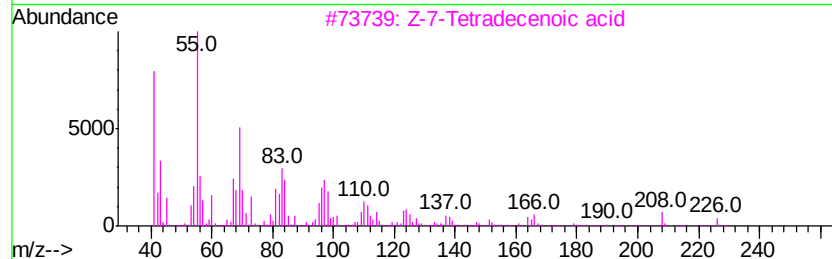
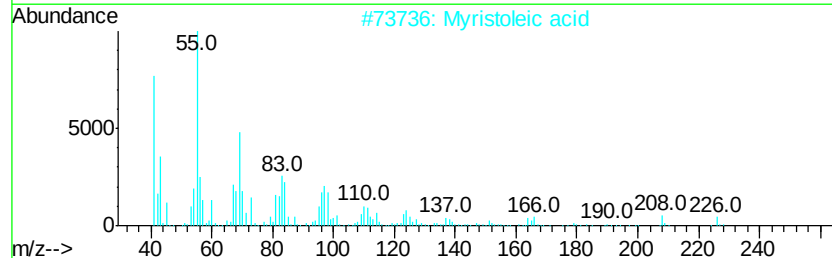
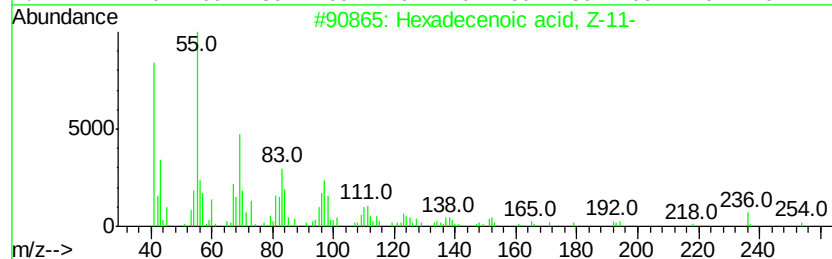
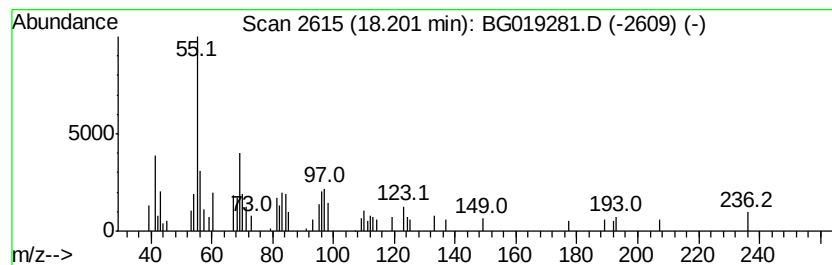
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 Peak Number 3 Hexadecenoic acid, Z-11- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.20	2.05 ng/ul	39675	Phenanthrene-d10	17.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	72
2		Myristoleic acid	226	C14H26O2	000544-64-9	49
3		Z-7-Tetradecenoic acid	226	C14H26O2	1000130-98-4	47
4		Cyclododecanemethanol	198	C13H26O	001892-12-2	43
5		Oxacyclotetradecane-2,11-dione, ...	240	C14H24O3	074685-36-2	43



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ALS Vial : 21 Sample Multiplier: 1

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BNA_G
ClientSampleId :
JG9K1

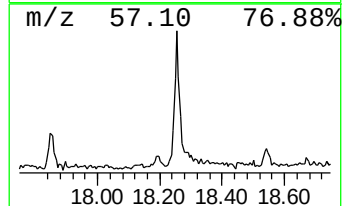
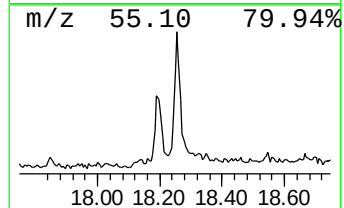
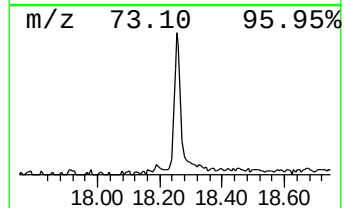
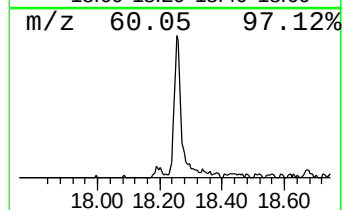
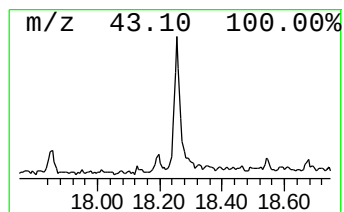
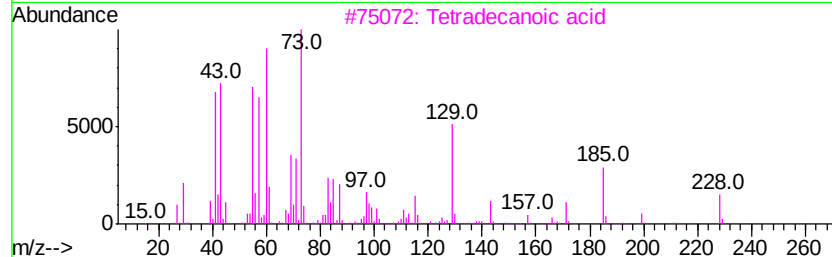
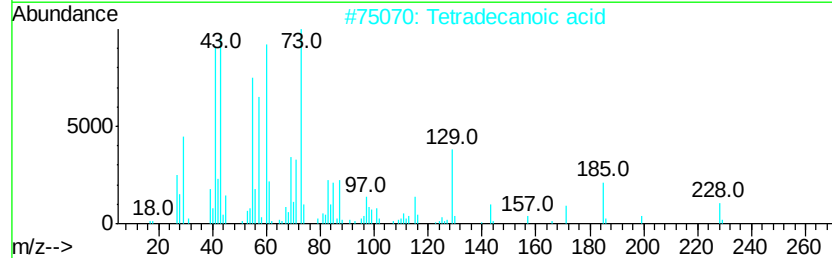
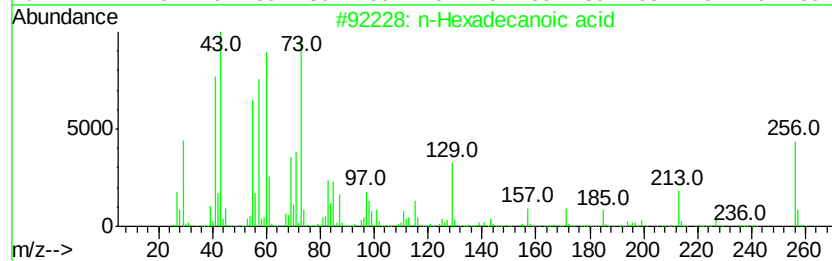
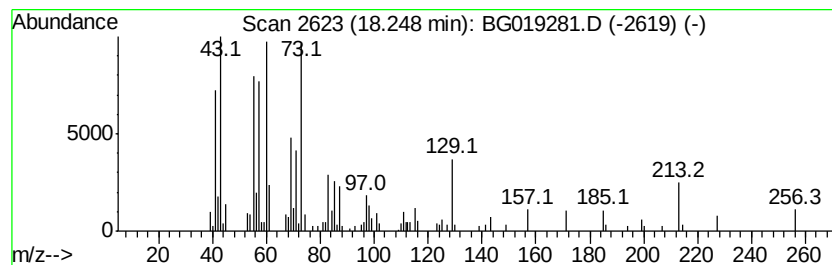
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.25	7.48 ng/ul	144615	Phenanthrene-d10	17.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	76
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	64



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
Data File : BG019281.D
Acq On : 21 Oct 2015 14:52
Operator : UM/NP
Sample : G4074-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9K1

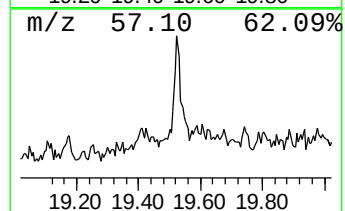
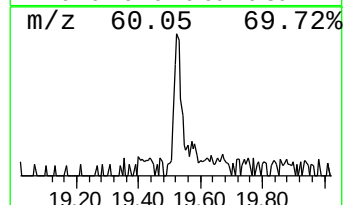
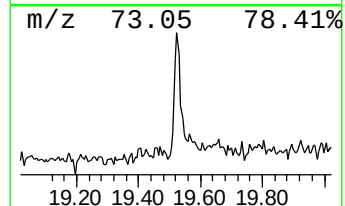
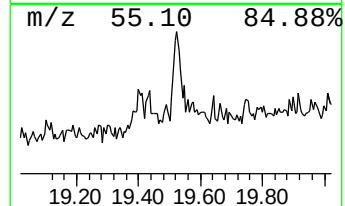
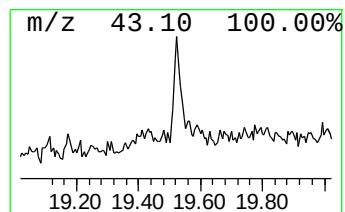
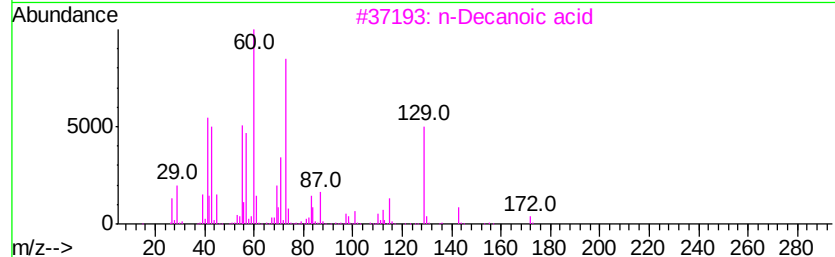
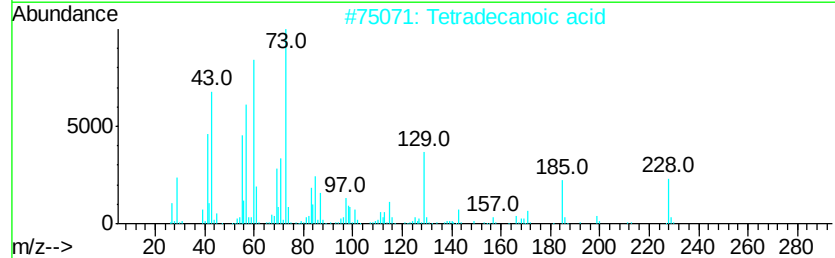
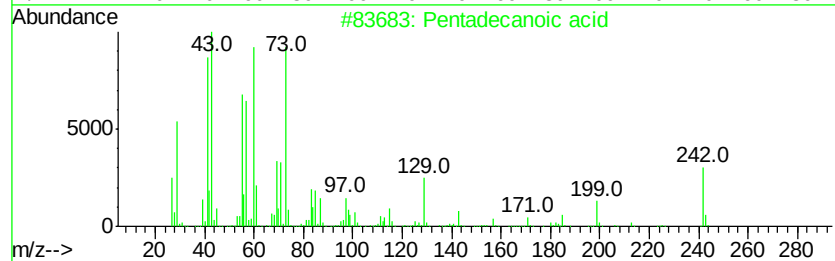
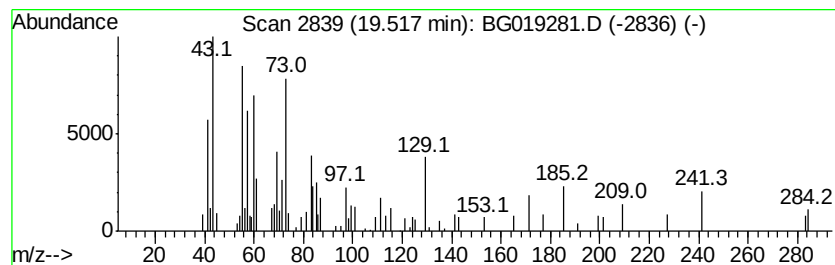
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.52	2.26 ng/ul	50849	Chrysene-d12	21.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecanoic acid	242	C15H30O2	001002-84-2	38
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	38
3		n-Decanoic acid	172	C10H20O2	000334-48-5	27
4		n-Decanoic acid	172	C10H20O2	000334-48-5	27
5		Undecanoic acid	186	C11H22O2	000112-37-8	27



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
Data File : BG019281.D
Acq On : 21 Oct 2015 14:52
Operator : UM/NP
Sample : G4074-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9K1

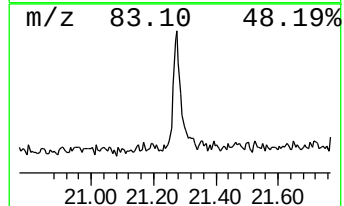
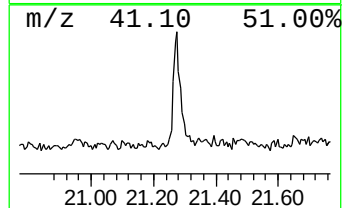
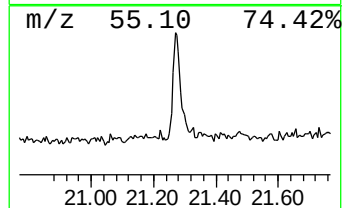
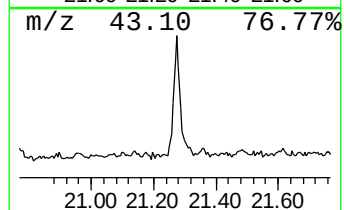
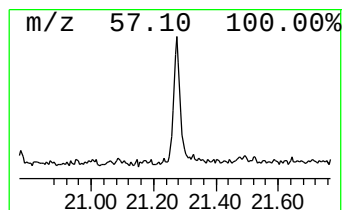
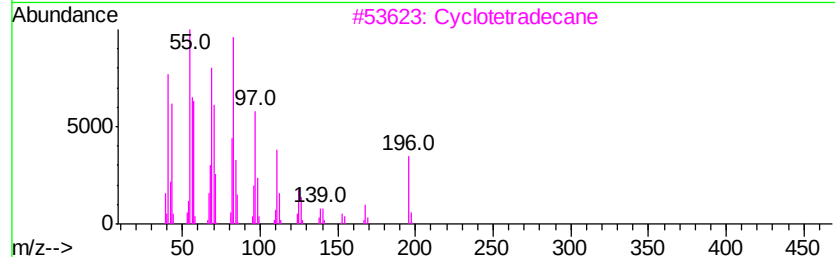
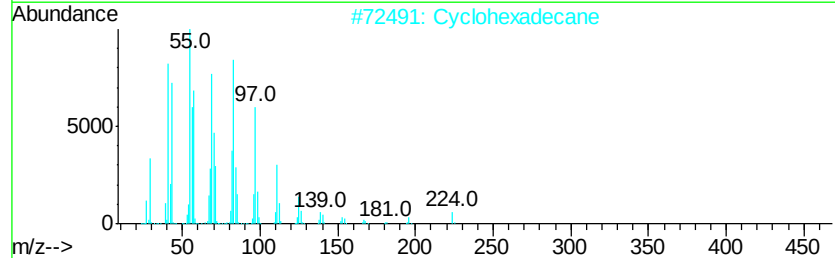
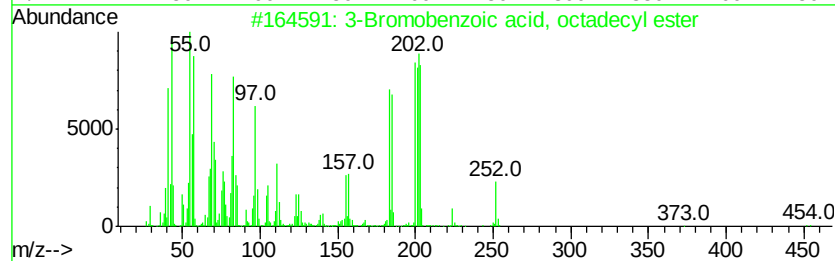
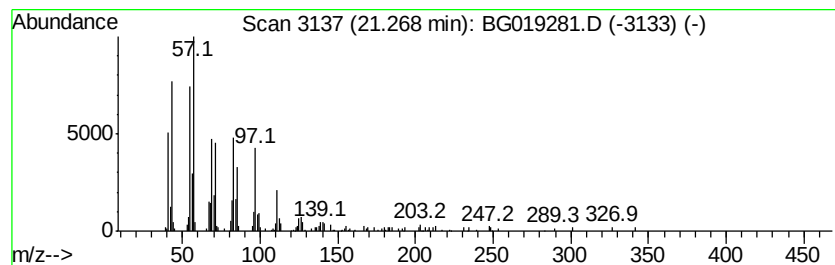
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 3-Bromobenzoic acid, octade... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.27	6.51 ng/ul	146801	Chrysene-d12	21.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Bromobenzoic acid, octadecyl e...	452	C25H41BrO2	070153-14-9	90
2		Cyclohexadecane	224	C16H32	000295-65-8	89
3		Cyclotetradecane	196	C14H28	000295-17-0	83
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	74
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
Data File : BG019281.D
Acq On : 21 Oct 2015 14:52
Operator : UM/NP
Sample : G4074-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9K1

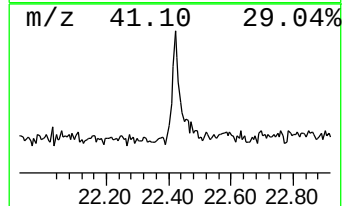
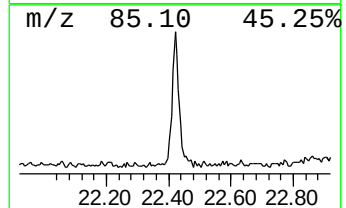
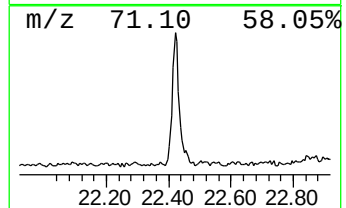
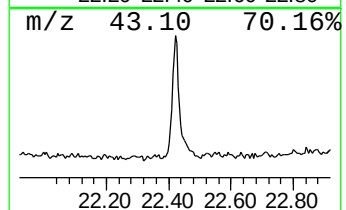
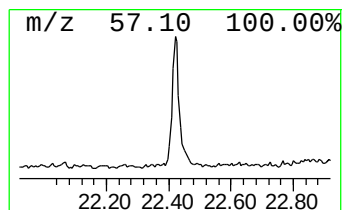
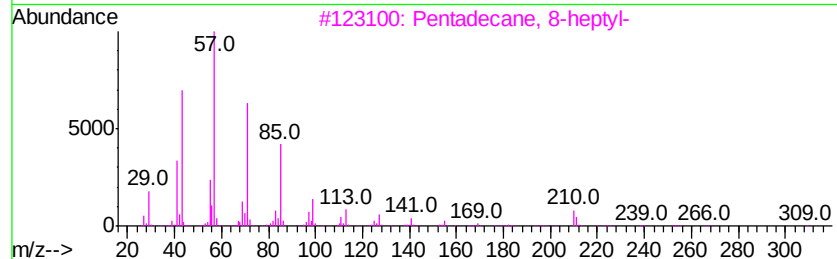
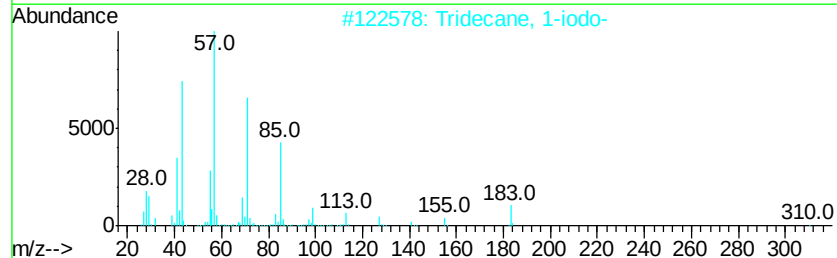
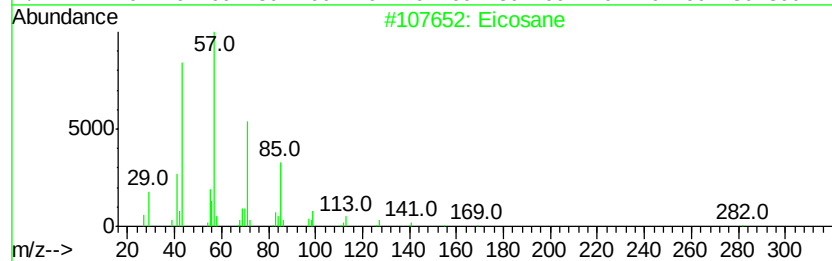
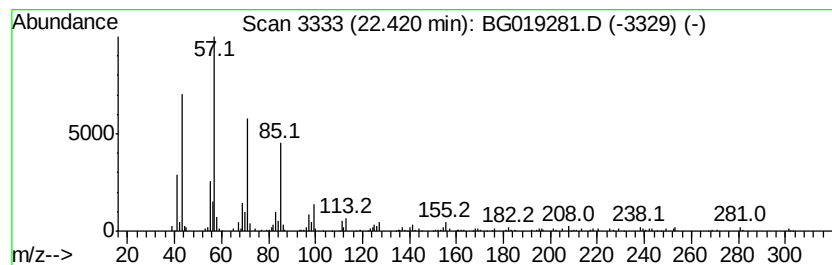
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.42	7.56 ng/ul	170376	Chrysene-d12	21.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
3			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83
4			Octacosane	394	C28H58	000630-02-4	83
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	83



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
Data File : BG019281.D
Acq On : 21 Oct 2015 14:52
Operator : UM/NP
Sample : G4074-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9K1

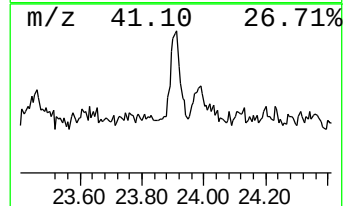
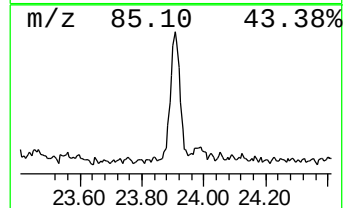
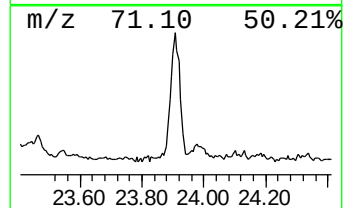
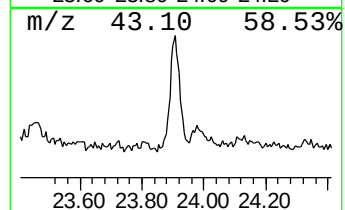
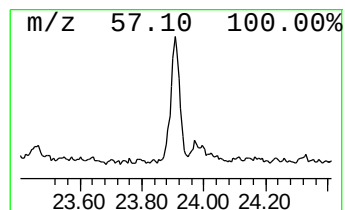
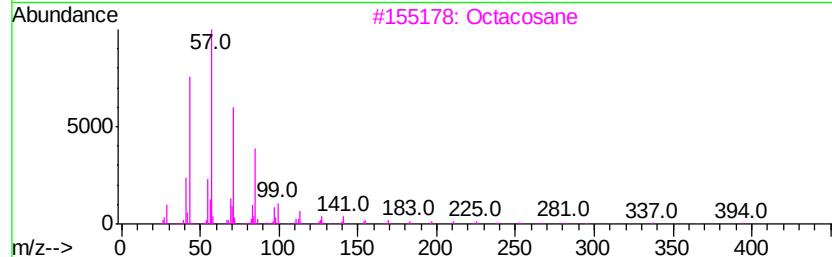
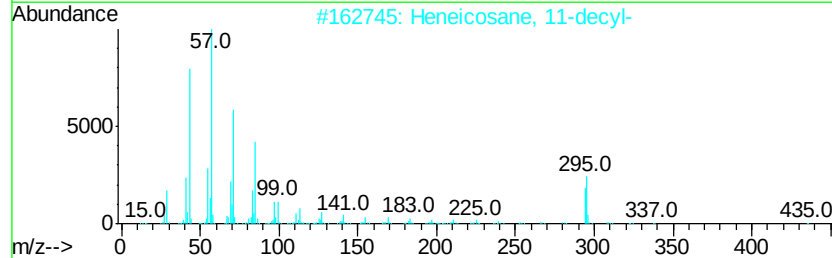
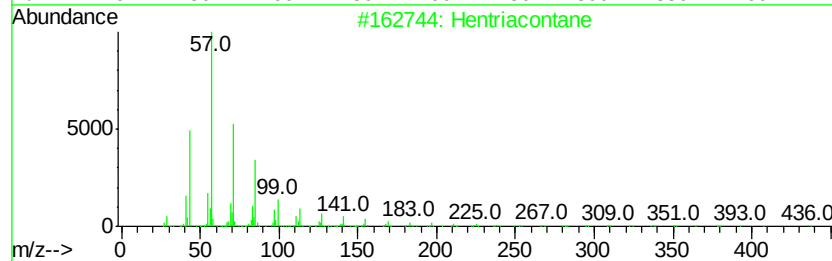
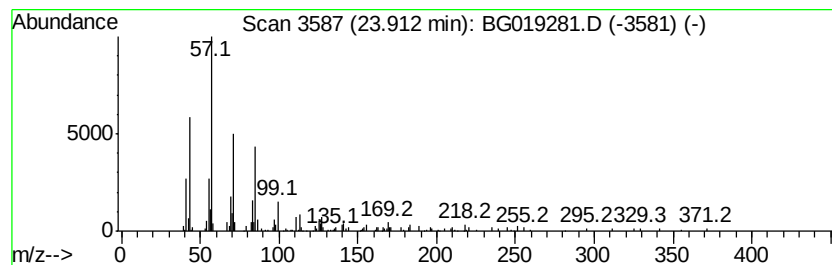
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.91	5.25 ng/ul	122149	Perylene-d12	24.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	87
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Hexacosane	366	C26H54	000630-01-3	86
5		Eicosane, 2-methyl-	296	C21H44	001560-84-5	86



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102115\
Data File : BG019281.D
Acq On : 21 Oct 2015 14:52
Operator : UM/NP
Sample : G4074-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9K1

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	3.08	4.6	ng/ul	27336	1	7.98	118452	20.0
Butane, 2-methoxy...	3.15	7.7	ng/ul	45381	1	7.98	118452	20.0
Hexadecenoic acid...	18.20	2.0	ng/ul	39675	4	17.37	386681	20.0
n-Hexadecanoic acid	18.25	7.5	ng/ul	144615	4	17.37	386681	20.0
unknown-01	19.52	2.3	ng/ul	50849	5	21.63	450828	20.0
3-Bromobenzoic ac...	21.27	6.5	ng/ul	146801	5	21.63	450828	20.0
(DEL) Alkane: Str...	22.42	7.6	ng/ul	170376	5	21.63	450828	20.0
(DEL) Alkane: Str...	23.91	5.3	ng/ul	122149	6	24.82	465347	20.0