

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
 Data File : BG024991.D
 Acq On : 6 Dec 2016 22:23
 Operator : UM/SJ
 Sample : H5843-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AH1

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 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: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.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	4.309	245	250	254	rBV5	8106	11225	0.27%	0.033%
2	4.362	254	259	268	rVB8	8932	22792	0.55%	0.067%
3	4.585	295	297	301	rBV3	8168	8133	0.19%	0.024%
4	5.313	417	421	425	rBV6	6238	8505	0.20%	0.025%
5	5.860	509	514	520	rVB8	6165	14945	0.36%	0.044%
6	5.919	520	524	528	rBV5	10085	13803	0.33%	0.041%
7	6.535	621	629	631	rBV6	5226	12434	0.30%	0.037%
8	6.841	677	681	685	rVB5	7111	9641	0.23%	0.028%
9	6.976	703	704	711	rVB4	5406	10035	0.24%	0.029%
10	7.217	738	745	748	rVV6	5756	10481	0.25%	0.031%
11	7.252	748	751	755	rVB3	5419	8627	0.21%	0.025%
12	7.329	758	764	766	rVB3	5622	8333	0.20%	0.024%
13	7.387	766	774	786	rBV4	50061	113026	2.71%	0.332%
14	7.558	796	803	814	rVB	183991	314194	7.52%	0.923%
15	7.769	832	839	848	rBV	193200	345211	8.26%	1.014%
16	8.145	899	903	907	rBV5	4899	8412	0.20%	0.025%
17	8.181	907	909	913	rVV4	7706	9238	0.22%	0.027%
18	8.245	913	920	932	rVB2	307310	568307	13.60%	1.669%
19	8.357	937	939	944	rVB4	5728	8217	0.20%	0.024%
20	8.410	944	948	951	rBV5	6877	9187	0.22%	0.027%
21	8.451	952	955	959	rVV6	7913	13215	0.32%	0.039%
22	8.662	988	991	996	rBV4	4872	8743	0.21%	0.026%
23	8.886	1026	1029	1031	rBV4	7056	7805	0.19%	0.023%
24	8.939	1031	1038	1049	rVB2	148440	289947	6.94%	0.852%
25	9.420	1109	1120	1132	rBV2	250537	482997	11.56%	1.419%
26	9.755	1170	1177	1180	rBV5	8946	13070	0.31%	0.038%
27	9.979	1213	1215	1220	rVB4	7045	8464	0.20%	0.025%
28	10.143	1235	1243	1252	rBV2	227857	418916	10.03%	1.230%
29	10.202	1252	1253	1258	rVB5	7413	7920	0.19%	0.023%
30	10.431	1287	1292	1294	rVB5	5216	8931	0.21%	0.026%
31	10.678	1325	1334	1349	rBV2	284018	585128	14.00%	1.719%
32	10.819	1357	1358	1362	rVB4	9331	8874	0.21%	0.026%
33	11.071	1390	1401	1413	rVB	409948	821215	19.65%	2.412%
34	11.160	1413	1416	1419	rBV2	7016	8333	0.20%	0.024%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	11.876	1536	1538	1543	rBV4	6368	9611	0.23%	0.028%
36	12.023	1556	1563	1568	rBV5	7074	14274	0.34%	0.042%
37	12.241	1594	1600	1606	rBV7	7868	17911	0.43%	0.053%
38	12.570	1654	1656	1664	rBV5	5555	11114	0.27%	0.033%
39	12.640	1664	1668	1671	rVV5	8407	14034	0.34%	0.041%
40	12.669	1671	1673	1676	rVB4	8585	8625	0.21%	0.025%
41	12.940	1716	1719	1723	rBV3	6049	9143	0.22%	0.027%
42	13.210	1761	1765	1768	rBV5	6966	10608	0.25%	0.031%
43	13.986	1895	1897	1906	rVB5	6723	16670	0.40%	0.049%
44	14.250	1928	1942	1949	rBV	752621	1177213	28.17%	3.458%
45	14.561	1980	1995	2004	rBV	742302	1202552	28.78%	3.532%
46	14.708	2015	2020	2027	rVB6	7417	15687	0.38%	0.046%
47	14.861	2039	2046	2056	rVB	834847	1320987	31.62%	3.880%
48	15.049	2070	2078	2090	rBV5	49046	129414	3.10%	0.380%
49	15.190	2101	2102	2105	rBV3	9139	7922	0.19%	0.023%
50	15.243	2108	2111	2117	rBV7	8570	16253	0.39%	0.048%
51	15.660	2176	2182	2185	rVB7	14773	25007	0.60%	0.073%
52	15.754	2195	2198	2203	rVB4	7906	11486	0.27%	0.034%
53	15.848	2203	2214	2226	rBV	1106429	1778047	42.55%	5.222%
54	15.960	2226	2233	2243	rVV	443274	719278	17.21%	2.113%
55	16.042	2243	2247	2250	rVV4	8525	15198	0.36%	0.045%
56	16.089	2250	2255	2260	rVB3	22543	32185	0.77%	0.095%
57	16.165	2261	2268	2270	rBV8	6751	8204	0.20%	0.024%
58	16.400	2305	2308	2310	rBV3	6904	7955	0.19%	0.023%
59	16.512	2324	2327	2334	rBV5	16958	31747	0.76%	0.093%
60	16.782	2366	2373	2374	rBV5	6236	12707	0.30%	0.037%
61	17.176	2437	2440	2447	rVB7	6927	13575	0.32%	0.040%
62	17.235	2447	2450	2454	rVB5	6978	10061	0.24%	0.030%
63	17.305	2455	2462	2466	rBV6	16504	25218	0.60%	0.074%
64	17.399	2468	2478	2486	rBV	2813994	4178297	100.00%	12.272%
65	17.464	2486	2489	2496	rVB8	12598	18321	0.44%	0.054%
66	17.546	2497	2503	2507	rBV3	137670	213577	5.11%	0.627%
67	17.605	2507	2513	2521	rVB2	1181244	1878298	44.95%	5.517%
68	17.699	2523	2529	2537	rBV2	1359894	2171471	51.97%	6.378%
69	17.758	2537	2539	2547	rVB8	17620	26249	0.63%	0.077%
70	17.934	2562	2569	2580	rBV2	70358	157392	3.77%	0.462%
71	18.040	2584	2587	2595	rVB8	9175	14284	0.34%	0.042%

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72	18.239	2612	2621	2626	rBV8	17614	32871	0.79%	0.097%
73	18.445	2651	2656	2661	rBV7	16168	24389	0.58%	0.072%
74	18.527	2665	2670	2673	rBV6	12778	20060	0.48%	0.059%
75	18.621	2685	2686	2690	rBV4	6739	8061	0.19%	0.024%
76	18.657	2690	2692	2697	rVB6	9094	13150	0.31%	0.039%
77	18.986	2742	2748	2749	rBV5	11775	17729	0.42%	0.052%
78	19.091	2764	2766	2769	rBV4	6894	9104	0.22%	0.027%
79	19.344	2806	2809	2811	rBV4	10465	9956	0.24%	0.029%
80	19.614	2849	2855	2859	rVB5	29616	46222	1.11%	0.136%
81	19.708	2867	2871	2877	rBV8	13398	30000	0.72%	0.088%
82	19.820	2885	2890	2904	rBV4	59988	180072	4.31%	0.529%
83	19.973	2910	2916	2926	rBV	1991112	2957634	70.79%	8.687%
84	20.167	2946	2949	2950	rBV3	8323	8076	0.19%	0.024%
85	21.729	3211	3215	3221	rVB5	39062	62049	1.49%	0.182%
86	21.894	3236	3243	3255	rBV	1575495	2719635	65.09%	7.988%
87	22.517	3345	3349	3356	rBV3	52149	130164	3.12%	0.382%
88	22.664	3370	3374	3381	rVB3	72405	115163	2.76%	0.338%
89	22.828	3398	3402	3405	rBV3	44277	70721	1.69%	0.208%
90	22.875	3406	3410	3417	rVB4	55457	87445	2.09%	0.257%
91	23.075	3438	3444	3454	rVB3	234174	547039	13.09%	1.607%
92	23.422	3492	3503	3513	rVB3	120602	366765	8.78%	1.077%
93	23.592	3526	3532	3546	rVB2	134044	292014	6.99%	0.858%
94	24.238	3635	3642	3648	rVB7	52647	113906	2.73%	0.335%
95	25.067	3772	3783	3796	rBV2	945668	2832611	67.79%	8.320%
96	25.237	3804	3812	3813	rBV2	81932	166449	3.98%	0.489%
97	25.302	3814	3823	3837	rVB2	876257	2685447	64.27%	7.887%
98	26.430	4005	4015	4024	rVB6	80482	268946	6.44%	0.790%
99	27.846	4249	4256	4275	rVB10	74587	320776	7.68%	0.942%
100	32.305	5004	5015	5032	rVB5	78588	381901	9.14%	1.122%

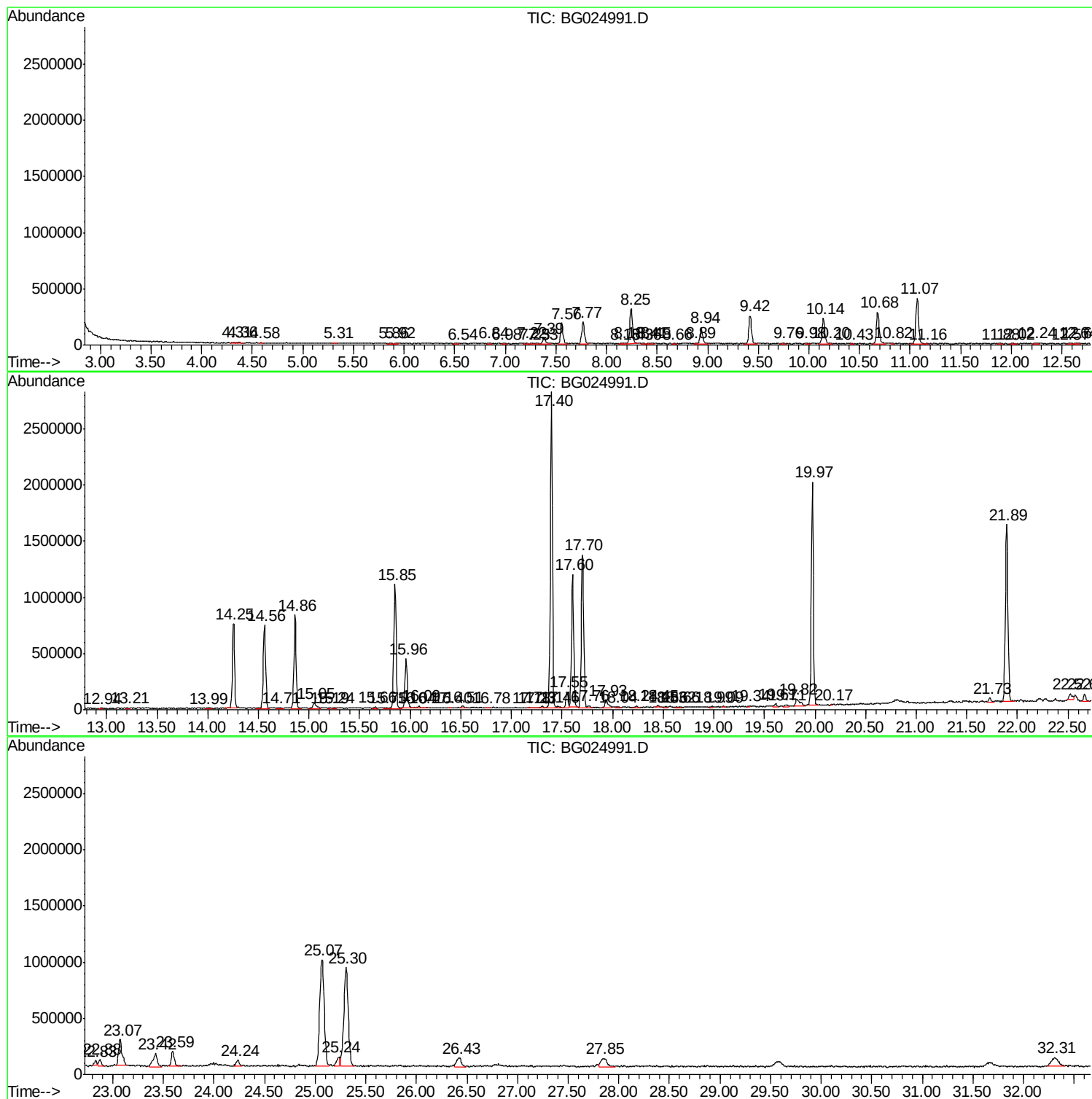
Sum of corrected areas: 34047234

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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



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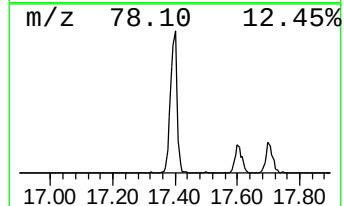
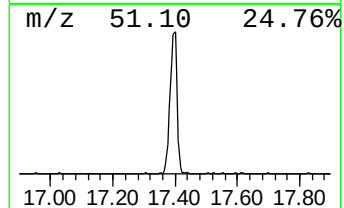
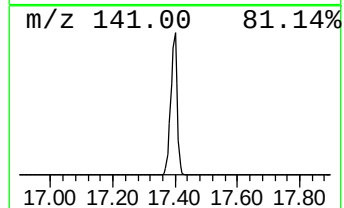
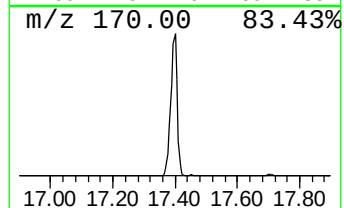
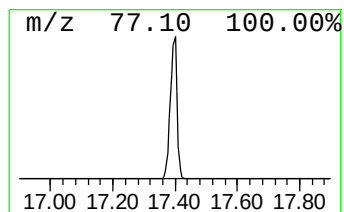
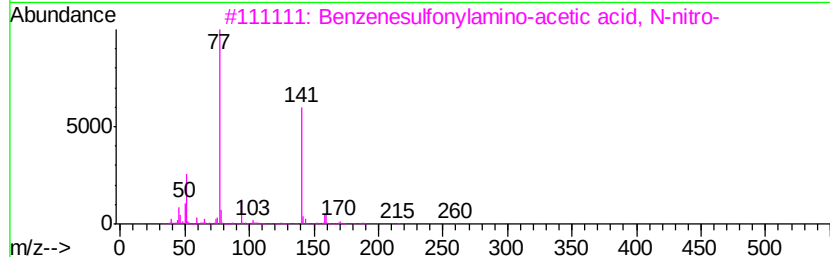
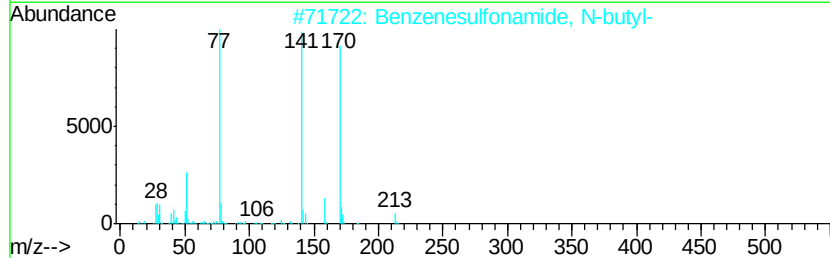
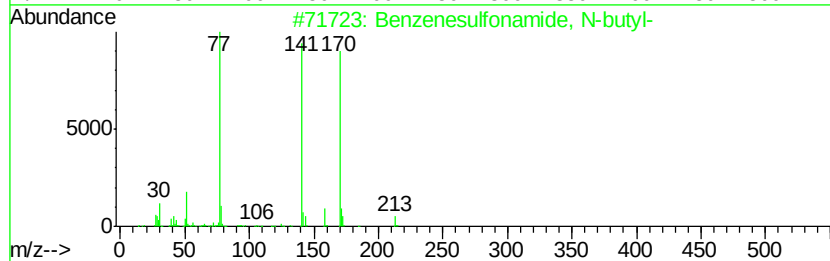
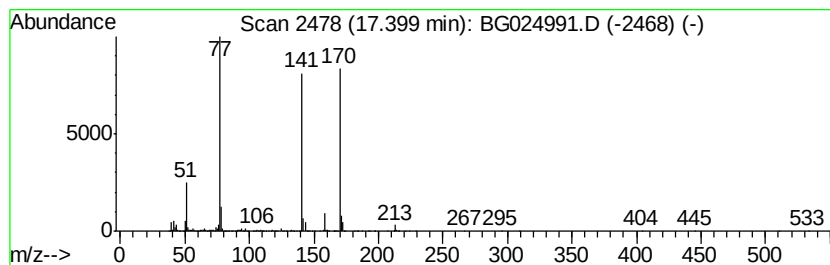
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzenesulfonamide, N-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.40	44.49 ng/ul	4178300	Phenanthrene-d10	17.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	90
2		Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	70
3		Benzenesulfonylamino-acetic acid...	260	C8H8N2O6S	1000301-09-0	53
4		N,N-Dichlorobenzenesulfonamide	225	C6H5Cl2NO2S	000473-29-0	50
5		N-Benzenesulfonyloxy-2,2-bis(tri...	335	C10H7F6NO3S	055734-41-3	50



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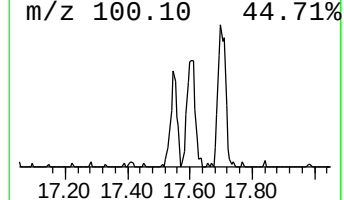
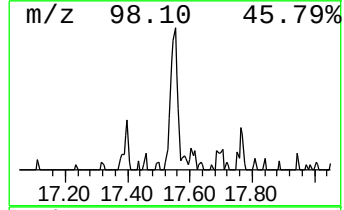
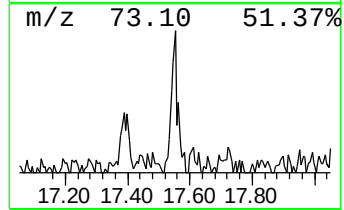
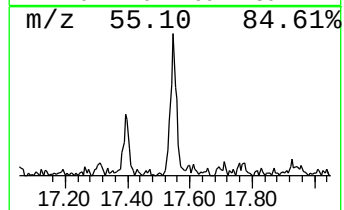
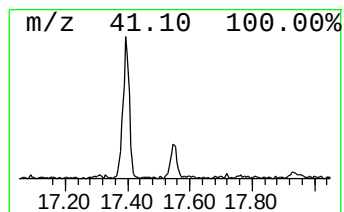
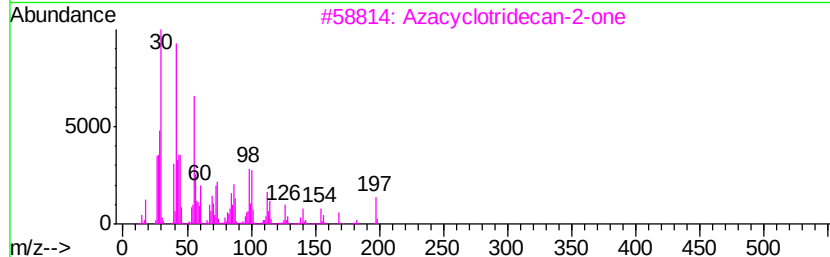
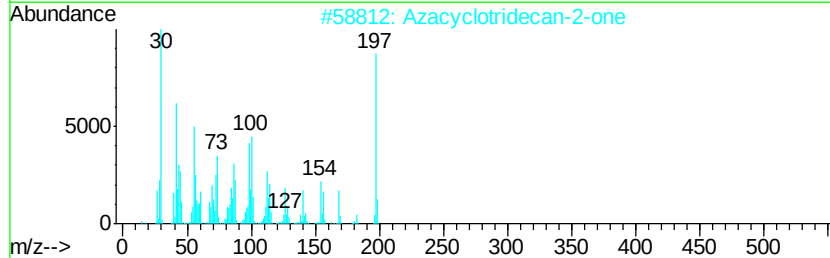
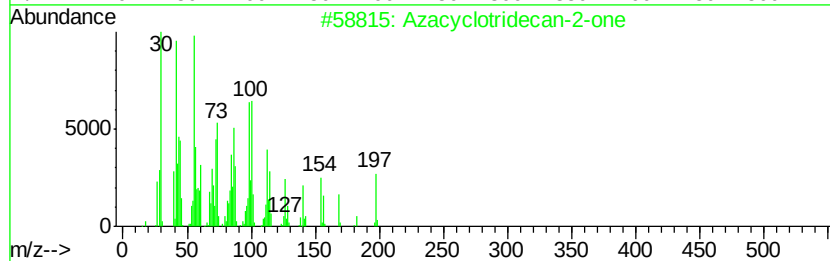
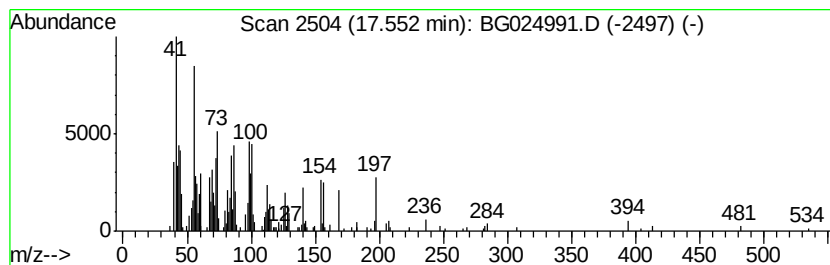
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Azacyclotridecan-2-one Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.55	2.27 ng/ul	213577	Phenanthrene-d10	17.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Azacyclotridecan-2-one	197	C12H23NO	000947-04-6	96
2		Azacyclotridecan-2-one	197	C12H23NO	000947-04-6	86
3		Azacyclotridecan-2-one	197	C12H23NO	000947-04-6	83
4		Azacyclotridecan-2-one	197	C12H23NO	000947-04-6	45
5		N-(3-Butyrylamino-propyl)-butyra...	214	C11H22N2O2	054535-62-5	35



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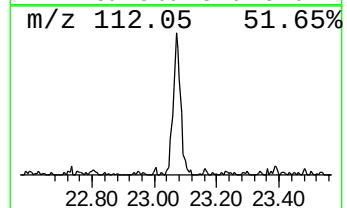
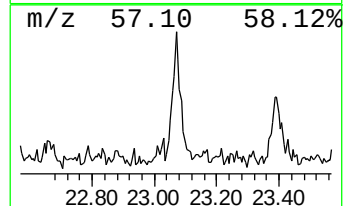
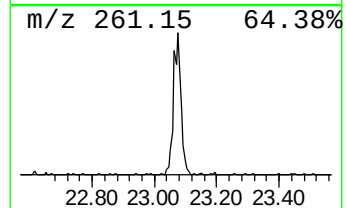
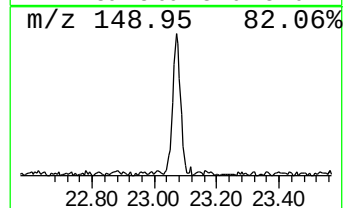
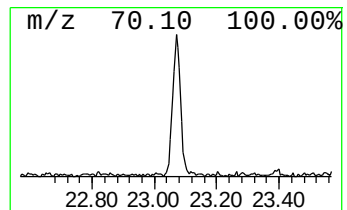
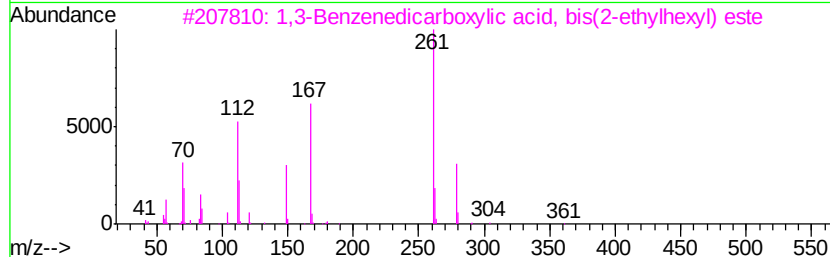
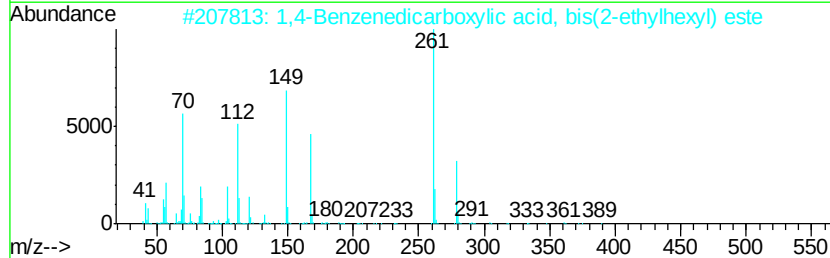
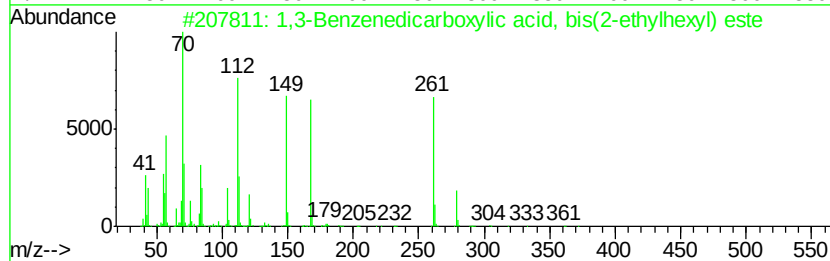
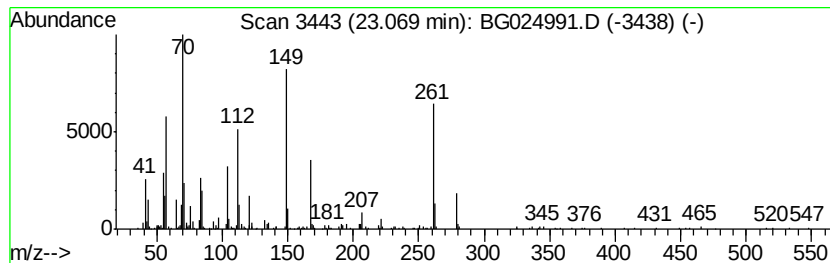
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Peak Number 3 1,3-Benzenedicarboxylic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.07	4.02 ng/ul	547039	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	80
2		1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	68
3		1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	47
4		1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	46
5		Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
Operator : UM/SJ
Sample : H5843-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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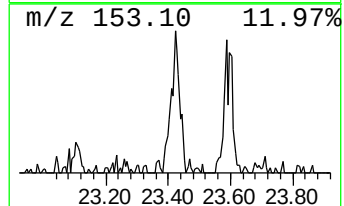
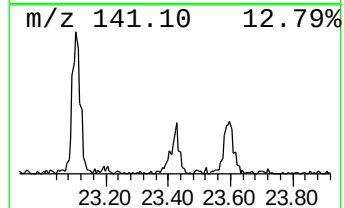
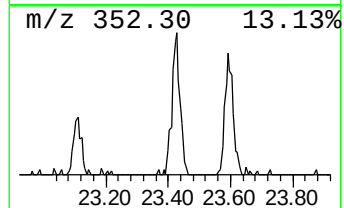
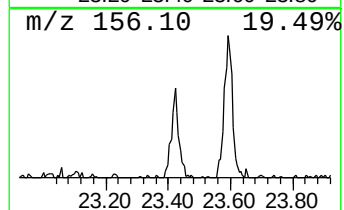
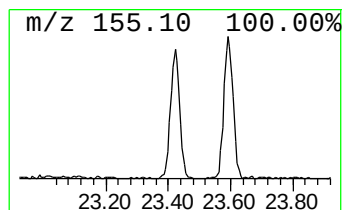
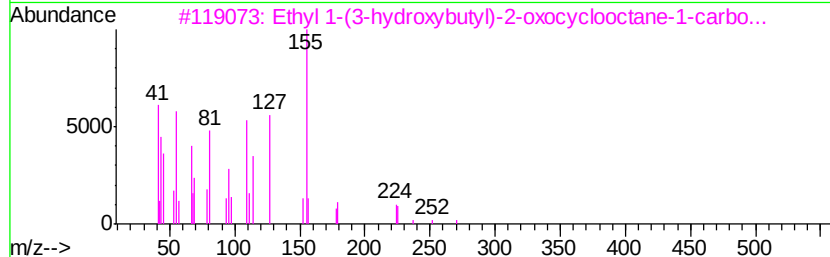
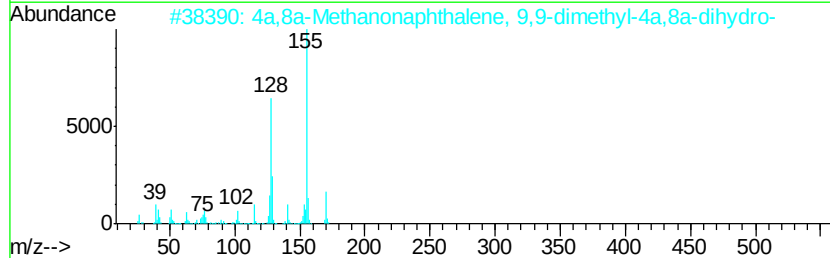
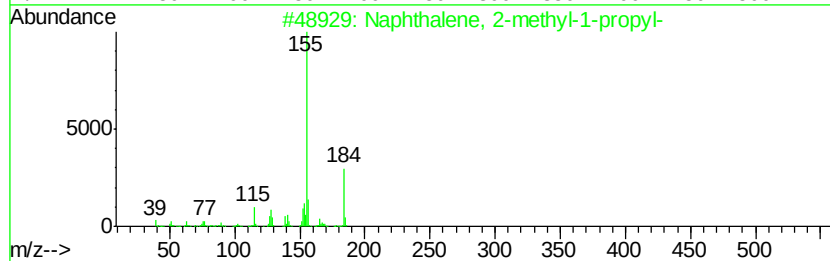
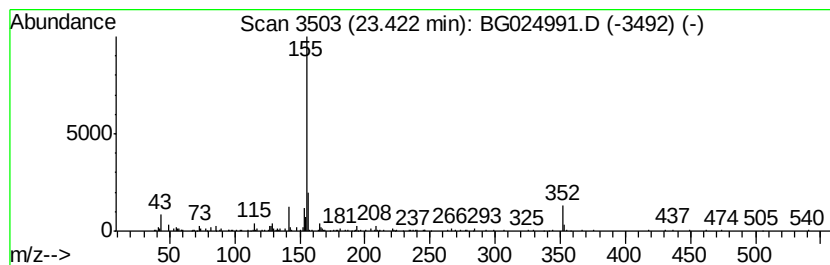
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2-methyl-1-pro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.42	2.70 ng/ul	366765	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-1-propyl-	184	C14H16	054774-89-9	64
2		4a,8a-Methanonaphthalene, 9,9-di...	170	C13H14	038963-97-2	40
3		Ethyl 1-(3-hydroxybutyl)-2-oxocyc...	270	C15H26O4	110288-16-9	38
4		1-Naphthoic acid, 3-chloroprop-2...	246	C14H11ClO2	1000308-82-8	36
5		1-Naphthoic acid, 2-biphenyl ester	324	C23H16O2	1000330-74-4	36



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
Operator : UM/SJ
Sample : H5843-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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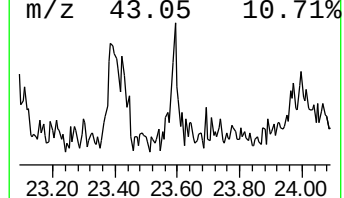
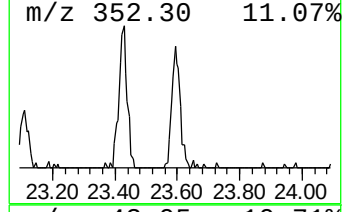
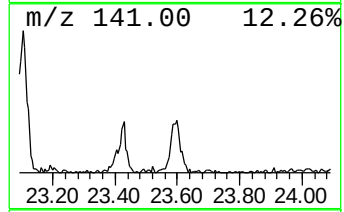
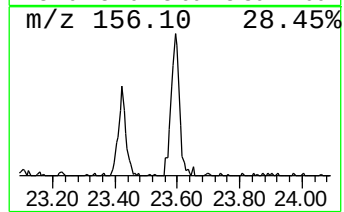
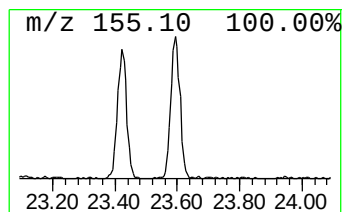
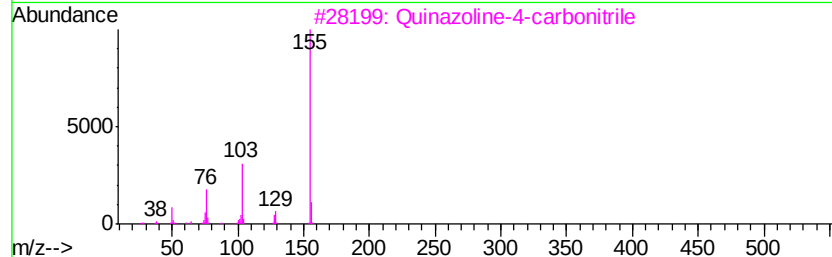
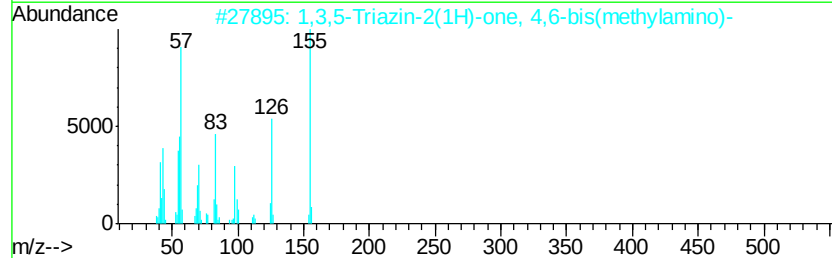
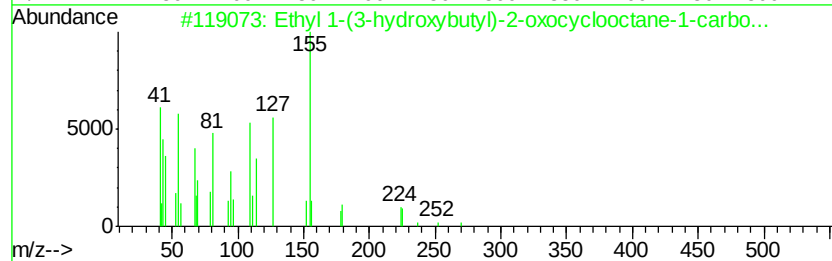
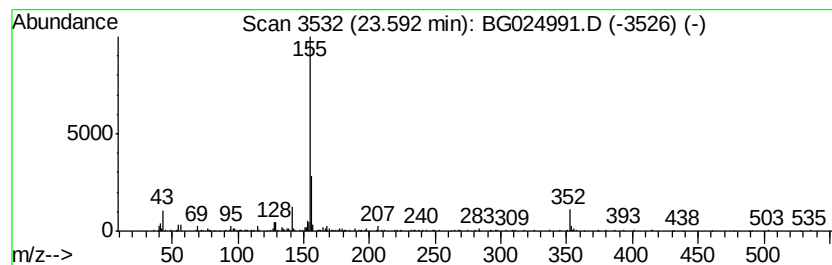
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethyl 1-(3-hydroxybutyl)-2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.59	2.15 ng/ul	292014	Chrysene-d12	21.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethyl 1-(3-hydroxybutyl)-2-oxocyclooctane-1-carboxylate	270	C15H26O4	110288-16-9	50
2		1,3,5-Triazin-2(1H)-one, 4,6-bis(methylamino)-	155	C5H9N5O	055702-52-8	38
3		Quinazoline-4-carbonitrile	155	C9H5N3	036082-71-0	38
4		1-Naphthoic acid, 2,4,5-trichloro-	350	C17H9Cl3O2	1000355-69-5	33
5		2,6-Difluoro-3-methylbenzoic acid	350	C14H7Cl3F2O2	1000357-29-9	28



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
Operator : UM/SJ
Sample : H5843-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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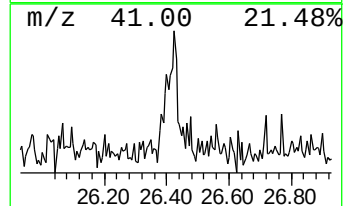
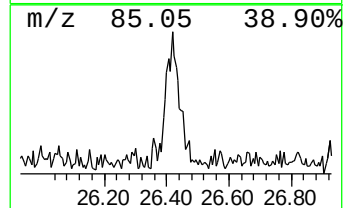
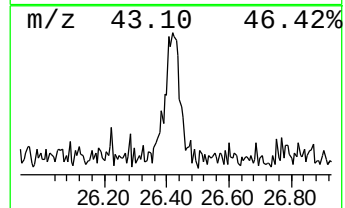
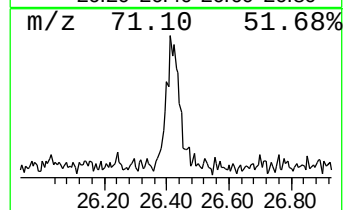
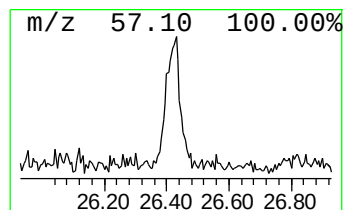
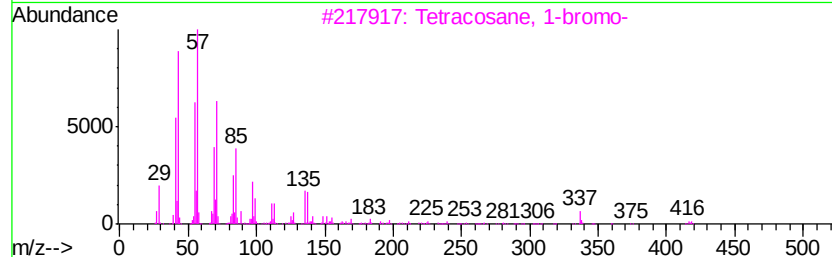
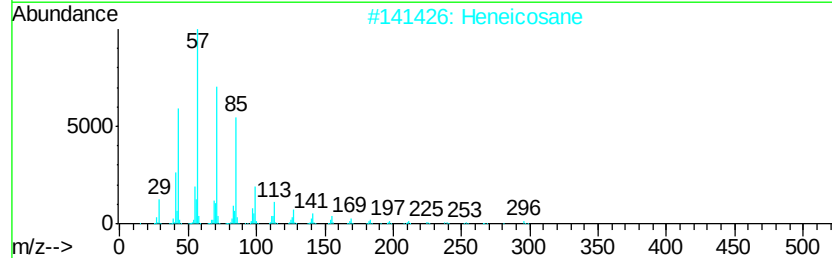
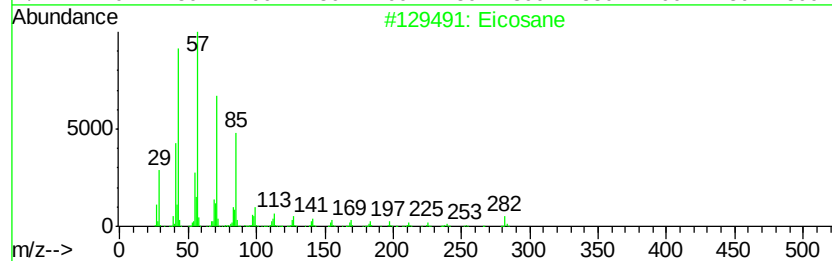
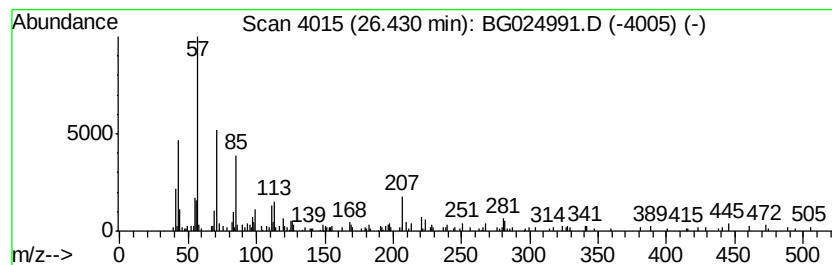
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.43	2.00 ng/ul	268946	Perylene-d12	25.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	89
2		Heneicosane	296	C21H44	000629-94-7	89
3		Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	68
4		7-Methyl-octadecane	268	C19H40	1000192-63-3	64
5		Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	64



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
Operator : UM/SJ
Sample : H5843-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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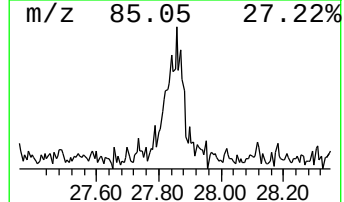
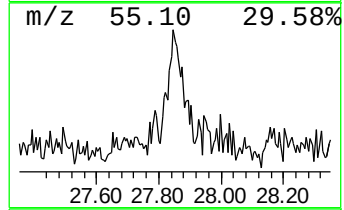
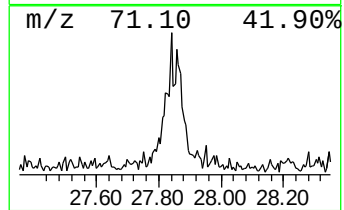
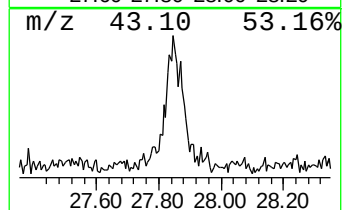
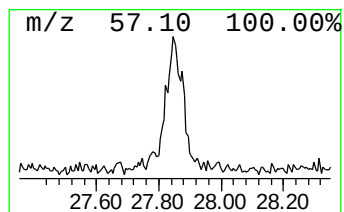
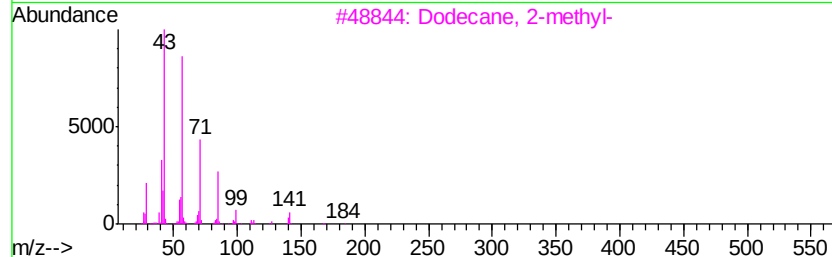
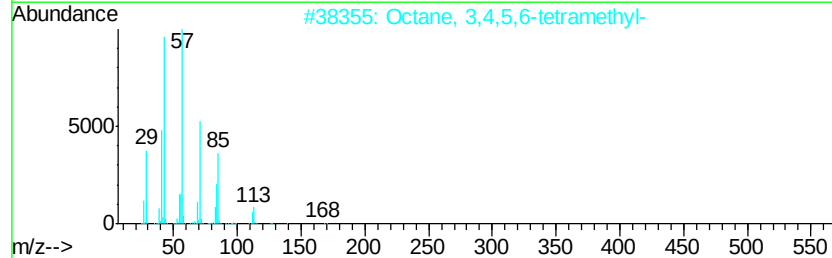
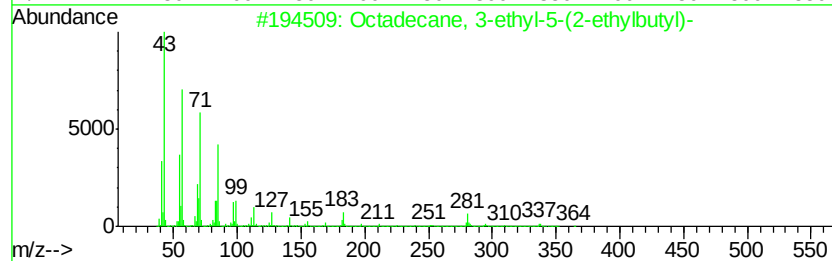
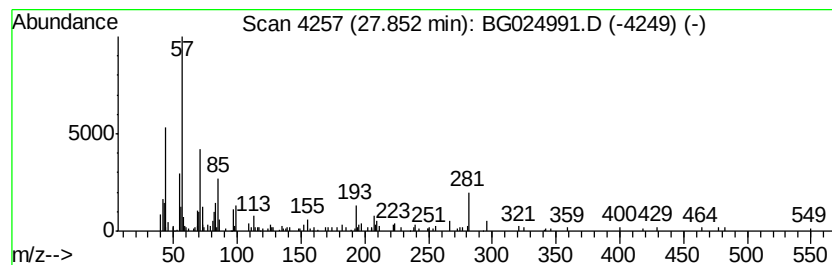
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.85	2.39 ng/ul	320776	Perylene-d12	25.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	27
2		Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	22
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	22
4		Piperazine, 1,4-diisobutyl-	226	C12H22N2O2	018940-58-4	22
5		Tetratetracontane	619	C44H90	007098-22-8	22



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
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Sample : H5843-02
Misc :
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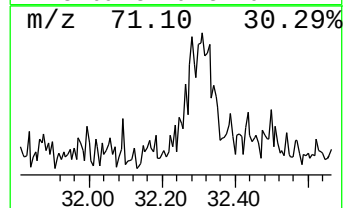
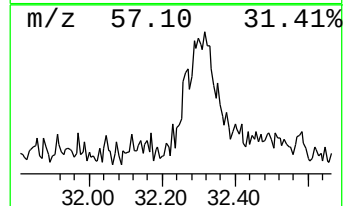
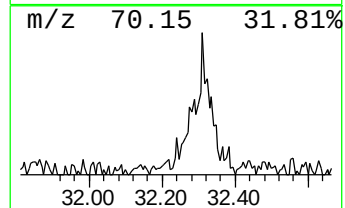
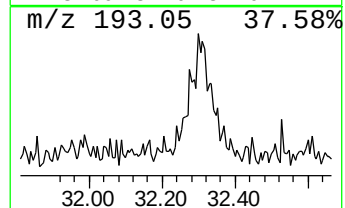
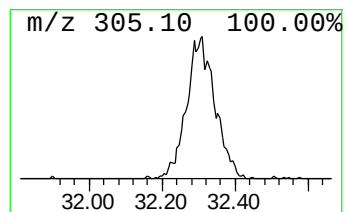
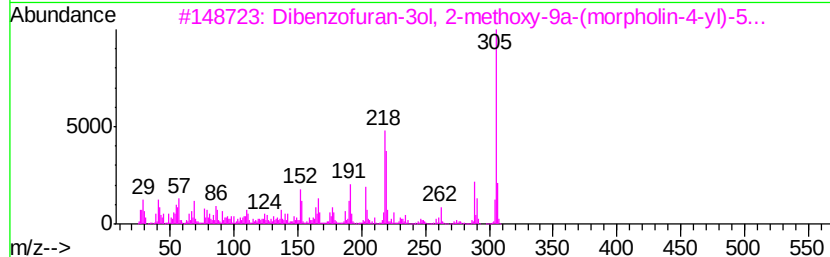
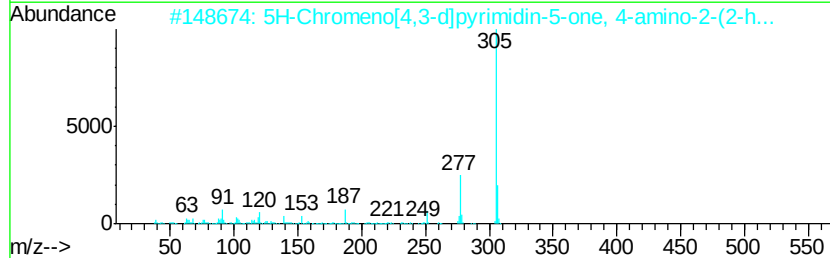
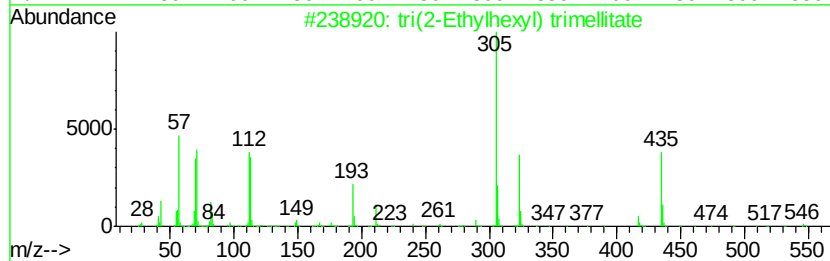
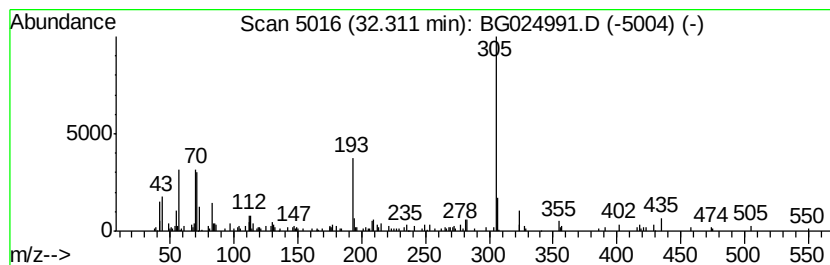
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 tri(2-Ethylhexyl) trimellitate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.		
32.31	2.84 ng/ul	381901	Perylene-d12		25.30		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	50
2			5H-Chromeno[4,3-d]pyrimidin-5-on...	305	C17H11N3O3	038370-06-8	49
3			Dibenzofuran-3-ol, 2-methoxy-9a-(...	305	C17H23NO4	094420-49-2	47
4			1-Diisopropyl octylsilylox-3,5-d...	348	C22H40OSi	1000307-85-1	33
5			tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	32



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120616\
Data File : BG024991.D
Acq On : 6 Dec 2016 22:23
Operator : UM/SJ
Sample : H5843-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Azacyclotridecan...	17.55	2.3	ng/ul	213577	4	17.60	1878300	20.0
1,3-Benzenedicarb...	23.07	4.0	ng/ul	547039	5	21.89	2719640	20.0
Naphthalene, 2-me...	23.42	2.7	ng/ul	366765	5	21.89	2719640	20.0
Ethyl 1-(3-hydrox...	23.59	2.1	ng/ul	292014	5	21.89	2719640	20.0
(DEL) Alkane: Str...	26.43	2.0	ng/ul	268946	6	25.30	2685450	20.0
(DEL) Alkane: Str...	27.85	2.4	ng/ul	320776	6	25.30	2685450	20.0
tri(2-Ethylhexyl)...	32.31	2.8	ng/ul	381901	6	25.30	2685450	20.0