

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
 Data File : BM001017.D
 Acq On : 16 Apr 2015 12:08
 Operator : TP/IZ
 Sample : G1749-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1N61

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_M\METHODS\SOM02.2-EPA-BM041515.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	2.787	13	17	27	rVB	1373650	1652099	17.79%	1.068%
2	2.946	41	44	49	rVB2	19923	20576	0.22%	0.013%
3	3.052	58	62	67	rVB3	24677	33654	0.36%	0.022%
4	4.151	246	249	253	rVB5	11987	14581	0.16%	0.009%
5	4.693	338	341	344	rBV5	11856	14387	0.15%	0.009%
6	5.804	524	530	534	rBV6	9047	15799	0.17%	0.010%
7	6.581	656	662	668	rVB	74349	112363	1.21%	0.073%
8	6.763	686	693	694	rBV	486709	675537	7.28%	0.437%
9	6.787	694	697	710	rVV	559378	894726	9.64%	0.578%
10	6.922	714	720	731	rVV	1001492	1525406	16.43%	0.986%
11	7.016	734	736	742	rVB6	8925	9446	0.10%	0.006%
12	7.116	745	753	755	rBV	1165762	1890809	20.36%	1.222%
13	7.145	755	758	764	rVV	1288673	1893764	20.40%	1.224%
14	7.210	764	769	777	rVB3	38584	76278	0.82%	0.049%
15	7.428	800	806	810	rVB2	16607	25048	0.27%	0.016%
16	7.581	820	832	843	rVB2	1148989	1902738	20.49%	1.230%
17	7.710	849	854	859	rVB	45126	71987	0.78%	0.047%
18	7.804	861	870	883	rVV	160871	304780	3.28%	0.197%
19	7.898	883	886	890	rVB3	8061	10903	0.12%	0.007%
20	7.998	899	903	909	rVB2	12324	17127	0.18%	0.011%
21	8.063	909	914	919	rBV2	32186	58747	0.63%	0.038%
22	8.116	919	923	927	rBV	93005	136539	1.47%	0.088%
23	8.151	927	929	936	rVB	35986	50316	0.54%	0.033%
24	8.251	940	946	949	rBV	96952	146816	1.58%	0.095%
25	8.298	949	954	962	rVB	1123975	1718911	18.51%	1.111%
26	8.392	962	970	978	rBV	2833782	4441146	47.83%	2.871%
27	8.457	978	981	984	rVB2	33378	41381	0.45%	0.027%
28	8.610	1002	1007	1015	rVB6	14597	29327	0.32%	0.019%
29	8.739	1022	1029	1037	rBV	1121662	1796758	19.35%	1.161%
30	8.857	1043	1049	1056	rVB	1307390	1997128	21.51%	1.291%
31	9.157	1094	1100	1105	rVB2	19285	33513	0.36%	0.022%
32	9.298	1117	1124	1135	rBV	1103358	1708310	18.40%	1.104%
33	9.457	1144	1151	1165	rVB	954244	1534356	16.53%	0.992%
34	9.616	1173	1178	1183	rBV5	10357	17066	0.18%	0.011%

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35	9.992	1235	1242	1255	rBV	1480418	2362106	25.44%	1.527%
36	10.363	1297	1305	1309	rBV	1442731	2474887	26.66%	1.600%
37	10.416	1309	1314	1321	rVV	2029026	3272581	35.25%	2.115%
38	10.510	1324	1330	1340	rVB	116771	204998	2.21%	0.133%
39	10.704	1356	1363	1371	rVB	2831761	4500996	48.48%	2.910%
40	11.663	1518	1526	1535	rBV	1646008	2590250	27.90%	1.674%
41	12.363	1639	1645	1651	rBV3	28024	41202	0.44%	0.027%
42	12.733	1701	1708	1722	rBV	998586	1535588	16.54%	0.993%
43	12.839	1722	1726	1729	rVV5	13756	22999	0.25%	0.015%
44	12.880	1729	1733	1739	rVB2	54413	75827	0.82%	0.049%
45	13.021	1754	1757	1762	rBV6	5649	9901	0.11%	0.006%
46	13.116	1768	1773	1781	rVB6	13651	27474	0.30%	0.018%
47	13.557	1843	1848	1854	rBV3	29329	40804	0.44%	0.026%
48	13.657	1859	1865	1872	rBV	2490328	3312122	35.67%	2.141%
49	13.880	1898	1903	1906	rBV2	201991	319786	3.44%	0.207%
50	13.933	1906	1912	1931	rVV	3017578	4425230	47.66%	2.861%
51	14.074	1933	1936	1941	rVB7	10102	14846	0.16%	0.010%
52	14.245	1958	1965	1978	rVB2	2013426	3043950	32.78%	1.968%
53	14.457	1995	2001	2011	rBV	336876	540149	5.82%	0.349%
54	14.568	2013	2020	2024	rVV	3001275	4207924	45.32%	2.720%
55	14.621	2024	2029	2031	rVV	1999438	2993308	32.24%	1.935%
56	14.645	2031	2033	2045	rVB	1552835	1774626	19.11%	1.147%
57	15.033	2096	2099	2104	rBV4	29138	34989	0.38%	0.023%
58	15.239	2128	2134	2141	rVV	3820983	5240772	56.45%	3.388%
59	15.368	2150	2156	2169	rVB	1247329	1664678	17.93%	1.076%
60	15.480	2169	2175	2180	rVB2	31750	44320	0.48%	0.029%
61	16.009	2258	2265	2272	rBV	53320	75146	0.81%	0.049%
62	16.092	2274	2279	2283	rVV5	8299	10471	0.11%	0.007%
63	16.304	2309	2315	2321	rBV	1863666	2567952	27.66%	1.660%
64	16.651	2367	2374	2387	rBV	2467318	3685851	39.70%	2.383%
65	16.774	2391	2395	2398	rVV5	14151	28532	0.31%	0.018%
66	16.809	2398	2401	2406	rVB2	22974	29108	0.31%	0.019%
67	16.974	2423	2429	2435	rBV2	4133144	7891026	84.99%	5.101%
68	17.033	2435	2439	2444	rVV	3260782	4389657	47.28%	2.838%
69	17.092	2444	2449	2452	rVV	3914109	5780202	62.25%	3.736%
70	17.121	2452	2454	2463	rVB	1860417	2233515	24.06%	1.444%
71	17.609	2533	2537	2543	rVB2	48269	62447	0.67%	0.040%

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72	17.856	2576	2579	2582	rVV4	9572	12710	0.14%	0.008%
73	17.898	2582	2586	2594	rVV6	15146	26947	0.29%	0.017%
74	18.168	2626	2632	2638	rBV	3151983	3965511	42.71%	2.563%
75	18.286	2643	2652	2657	rBV2	4815986	9284720	100.00%	6.002%
76	18.380	2665	2668	2671	rBV	17822	19018	0.20%	0.012%
77	18.415	2671	2674	2678	rVV6	10202	19031	0.20%	0.012%
78	18.462	2678	2682	2685	rVB5	8731	12090	0.13%	0.008%
79	19.056	2777	2783	2792	rBV	2925605	3669327	39.52%	2.372%
80	19.192	2802	2806	2809	rBV5	7431	13316	0.14%	0.009%
81	19.280	2815	2821	2822	rBV	34666	42510	0.46%	0.027%
82	19.392	2834	2840	2843	rBV2	4492940	6538692	70.42%	4.227%
83	19.421	2843	2845	2855	rVB	3245785	3699676	39.85%	2.392%
84	19.721	2891	2896	2904	rBV	4541380	5099731	54.93%	3.297%
85	19.809	2908	2911	2917	rVB6	7852	10368	0.11%	0.007%
86	19.897	2921	2926	2931	rBV	4190049	4816368	51.87%	3.113%
87	20.327	2995	2999	3003	rBV3	35021	42341	0.46%	0.027%
88	20.621	3046	3049	3052	rBV5	10654	11639	0.13%	0.008%
89	20.903	3094	3097	3103	rVB8	12268	20452	0.22%	0.013%
90	20.986	3107	3111	3115	rBV3	33409	40080	0.43%	0.026%
91	21.115	3128	3133	3140	rBV	4526720	4810625	51.81%	3.110%
92	21.192	3140	3146	3155	rVV	2857415	3566123	38.41%	2.305%
93	23.227	3485	3492	3507	rVB2	3204637	5746371	61.89%	3.715%
94	23.362	3508	3515	3524	rBV	1885162	3361643	36.21%	2.173%
95	25.544	3876	3886	3900	rBV	1373030	3466575	37.34%	2.241%

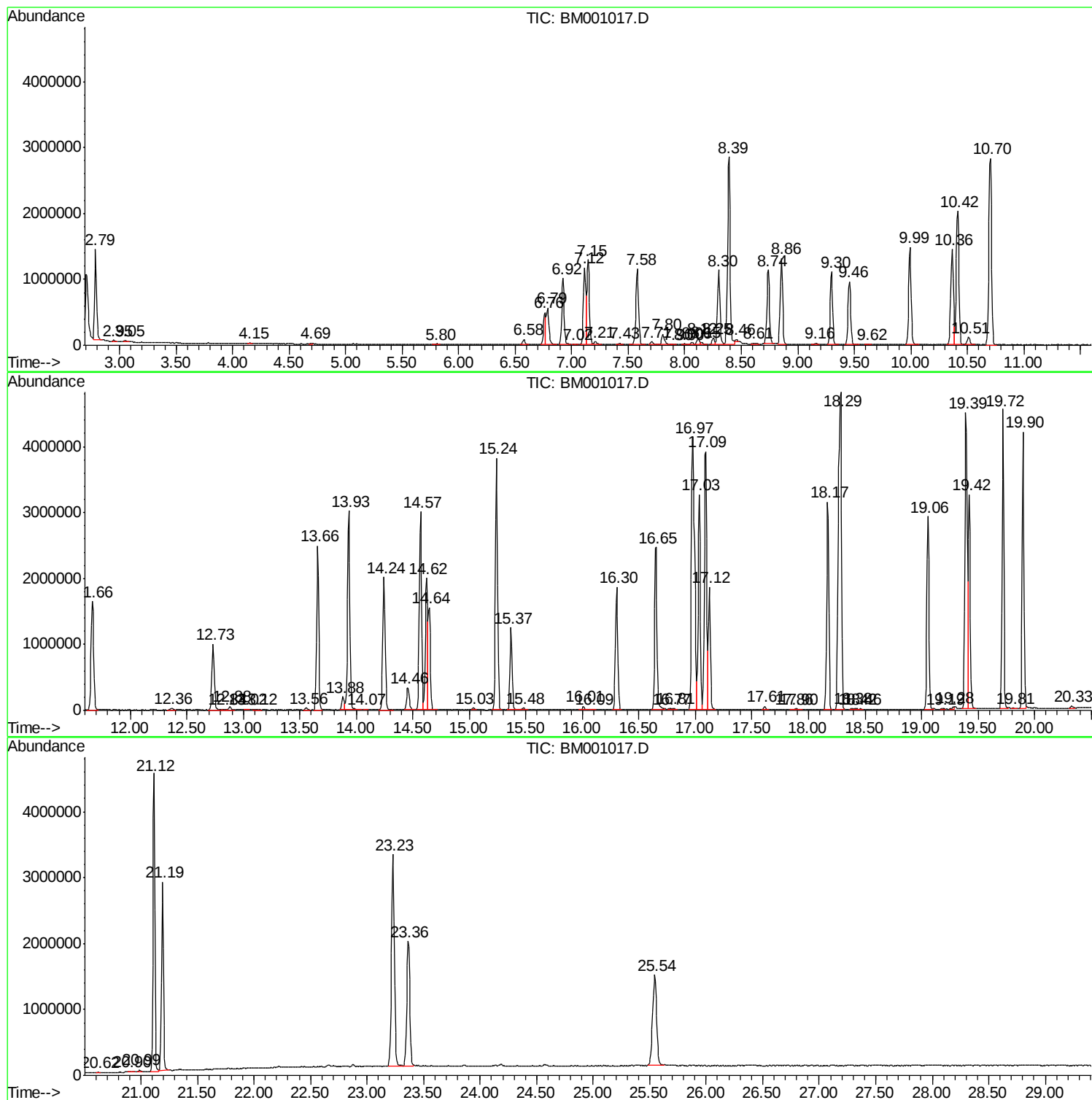
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM041515.M
Quant Title : SVOA CALIBRATION

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



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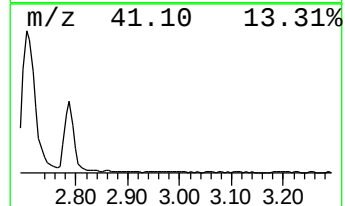
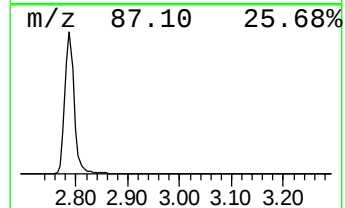
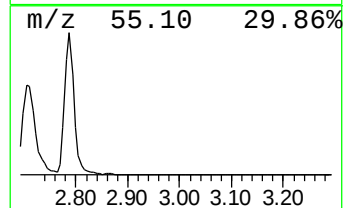
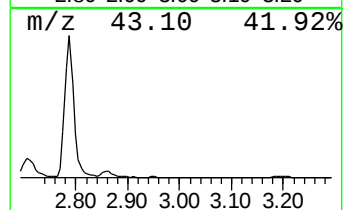
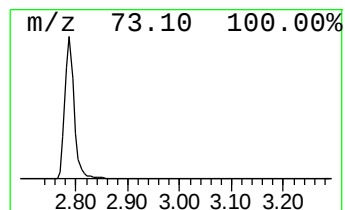
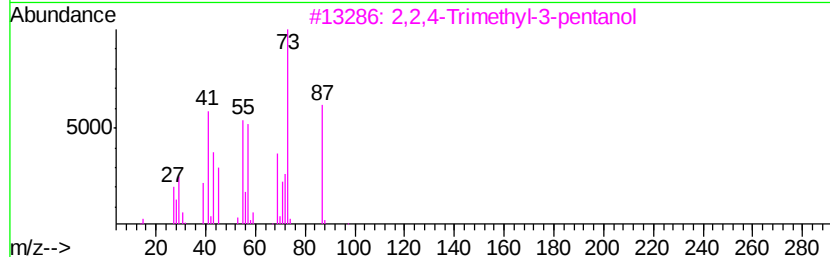
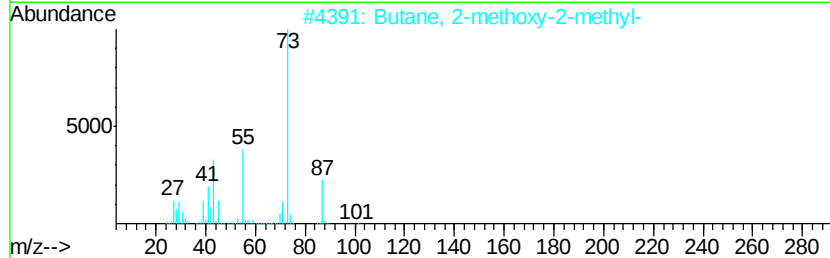
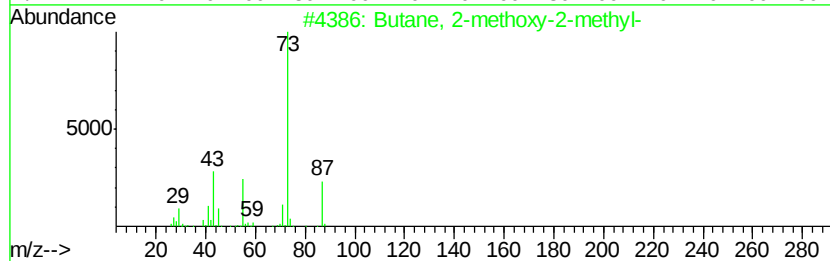
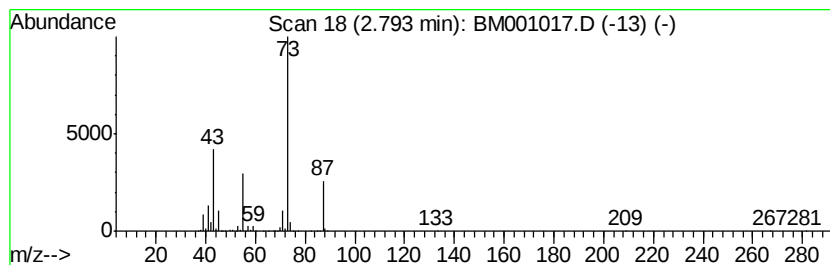
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM041515.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	17.37 ng/ul	1652100	1,4-Dichlorobenzene-d4	7.58

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		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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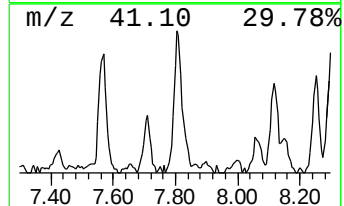
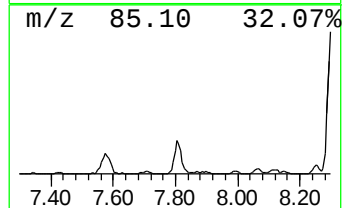
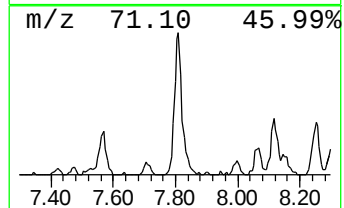
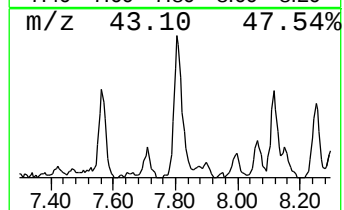
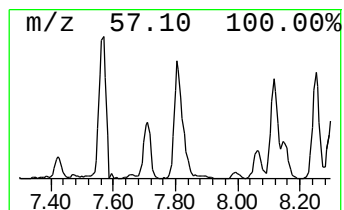
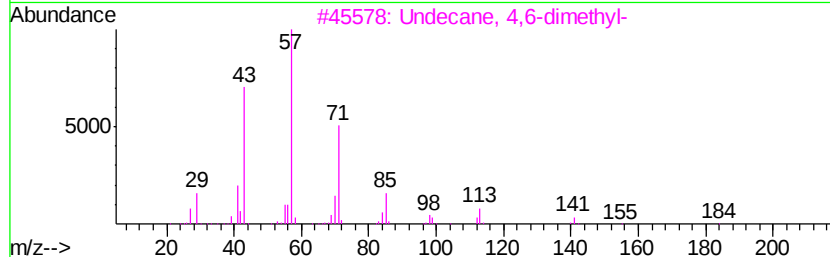
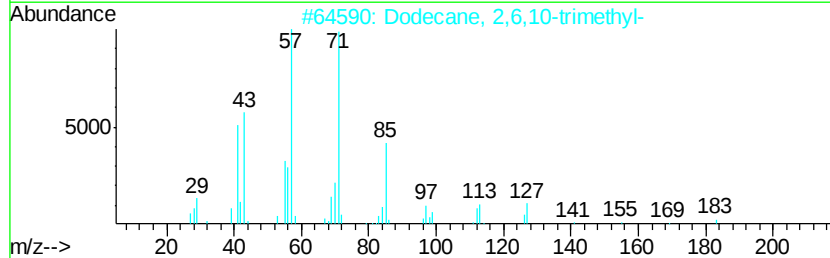
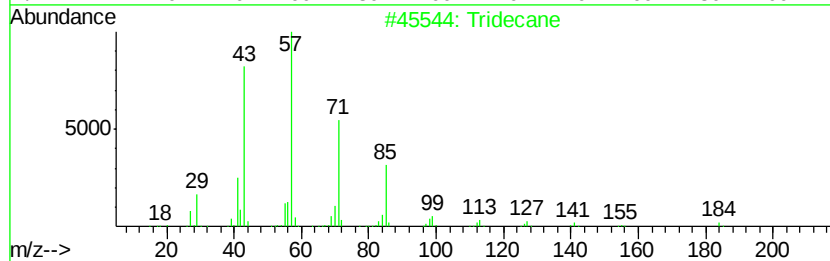
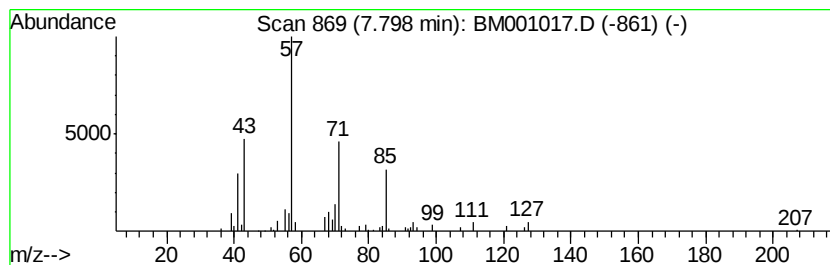
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chain Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	3.20 ng/ul	304780	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	64
2		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
3		Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	64
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
5		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	59



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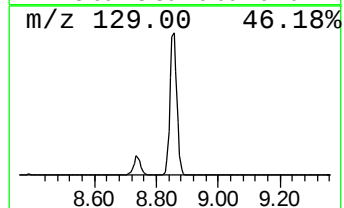
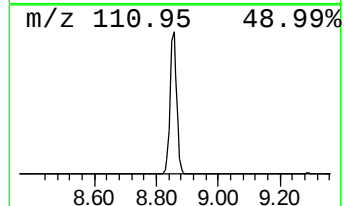
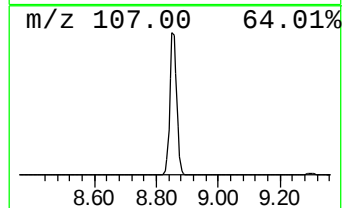
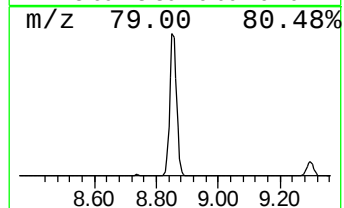
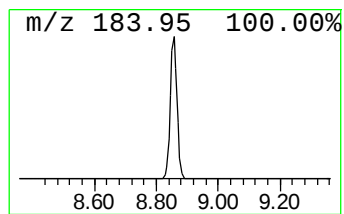
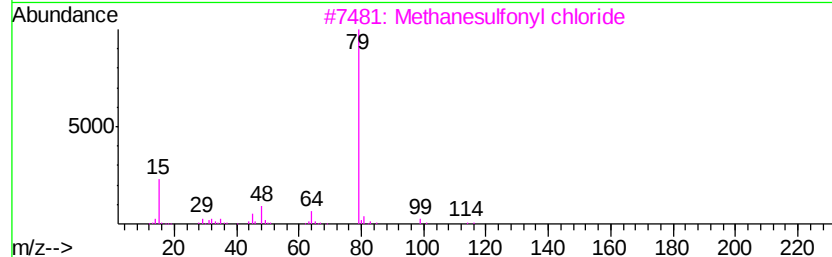
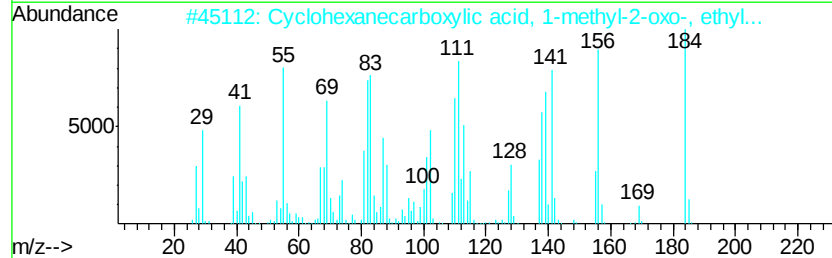
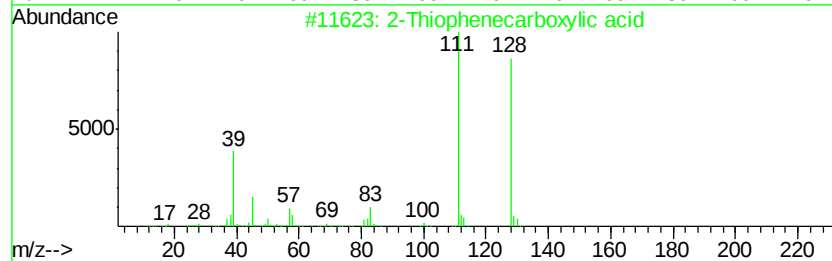
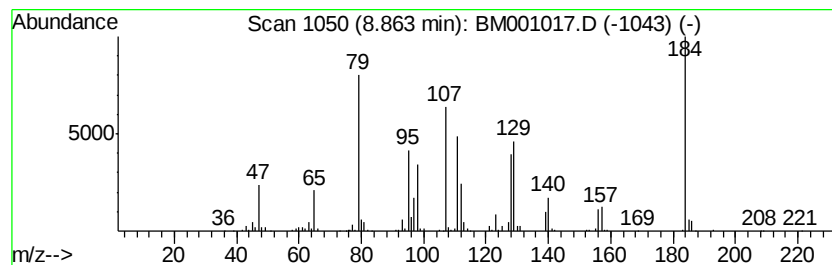
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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.86	20.99 ng/ul	1997130	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Thiophenecarboxylic acid	128	C5H4O2S	000527-72-0	11
2		Cyclohexanecarboxylic acid, 1-me...	184	C10H16O3	005453-94-1	10
3		Methanesulfonyl chloride	114	CH3ClO2S	000124-63-0	10
4		2-Oxa-7-thiatricyclo[4.4.0.0(3,8...	156	C8H12O5	039637-14-4	10
5		Dibenzothiophene	184	C12H8S	000132-65-0	10



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Misc :
ALS Vial : 4 Sample Multiplier: 1

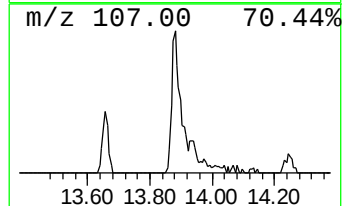
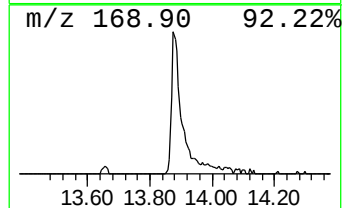
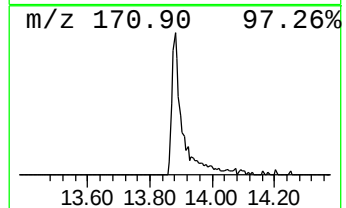
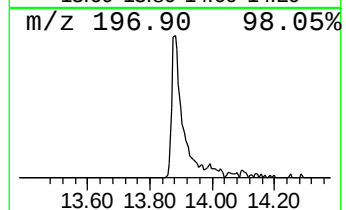
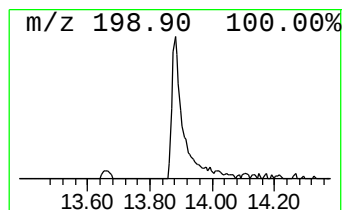
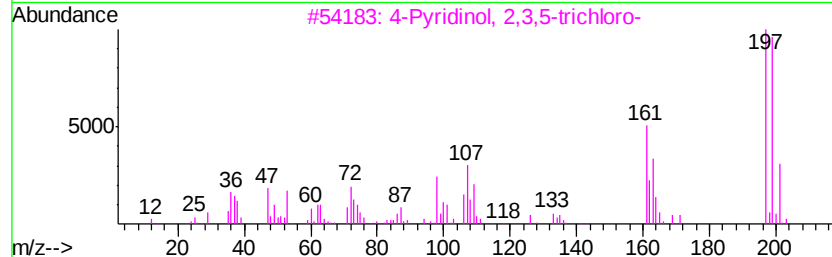
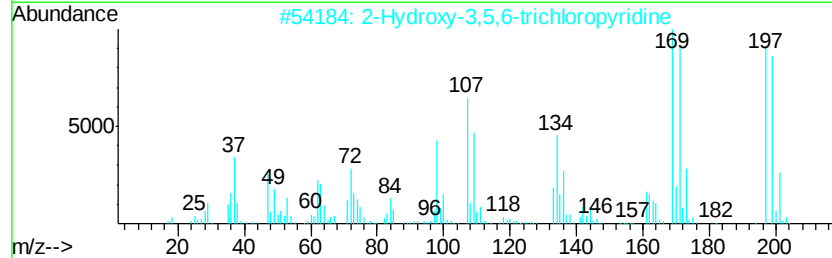
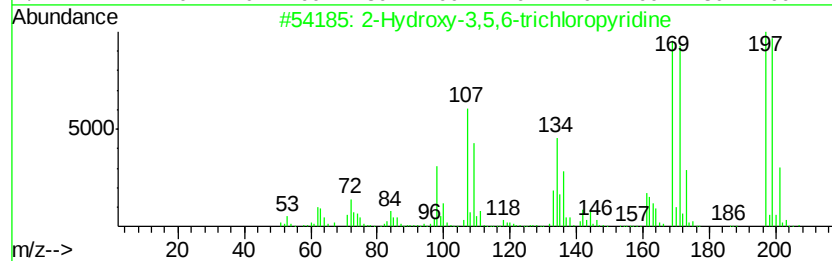
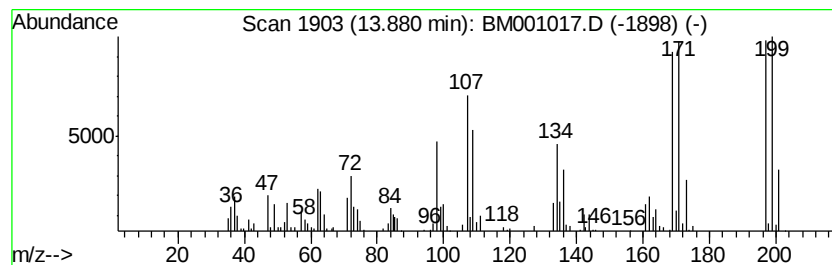
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM041515.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Hydroxy-3,5,6-trichloropy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	2.10 ng/ul	319786	Acenaphthene-d10	14.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		2-Hydroxy-3,5,6-trichloropyridine	197 C5H2Cl3NO	006515-38-4 99
2		2-Hydroxy-3,5,6-trichloropyridine	197 C5H2Cl3NO	006515-38-4 87
3		4-Pyridinol, 2,3,5-trichloro-	197 C5H2Cl3NO	001970-40-7 51
4		2-Propenenitrile, 3-(2,4-dichlor...	197 C9H5Cl2N	095727-84-7 38
5		Benzothiazole, 2-chloro-	169 C7H4ClNS	000615-20-3 12



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
Data File : BM001017.D
Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
X1N61

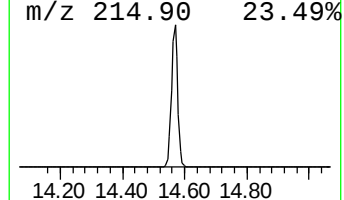
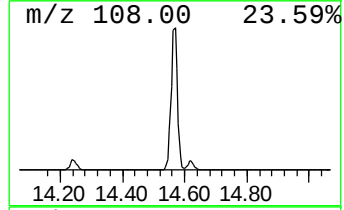
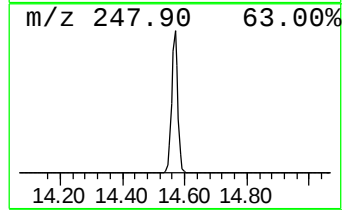
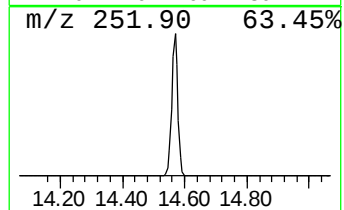
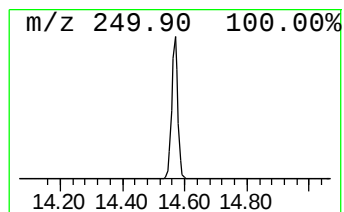
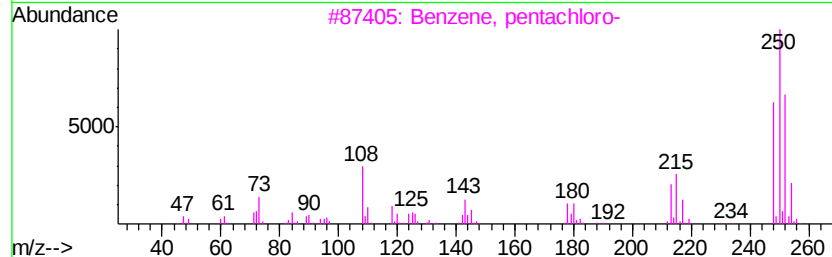
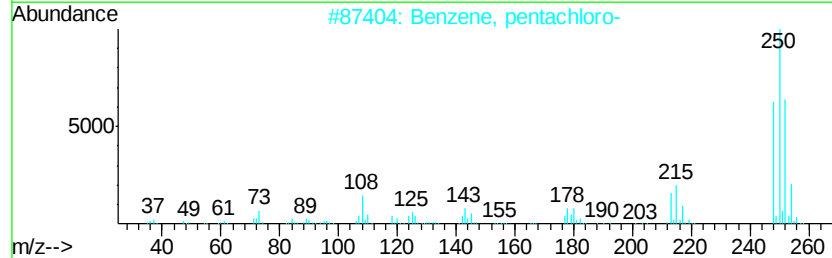
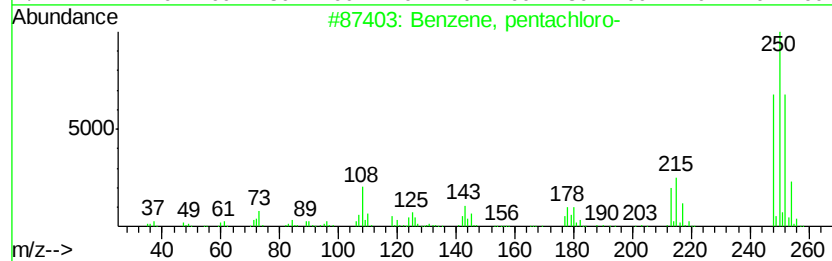
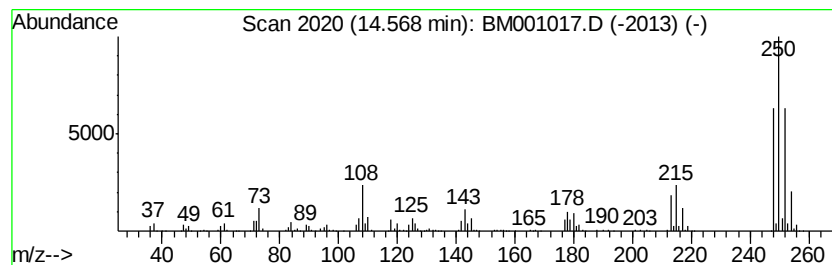
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, pentachloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	27.65 ng/ul	4207920	Acenaphthene-d10	14.24

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
2	Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
3	Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
4	Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
5	2,3-Dichloro-1,4,5,8-tetraaza-ph...	250	C10H4Cl2N4	071222-45-2	53



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
Data File : BM001017.D
Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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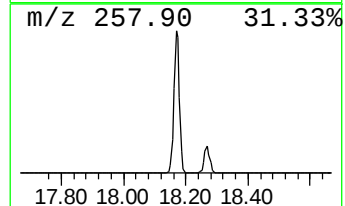
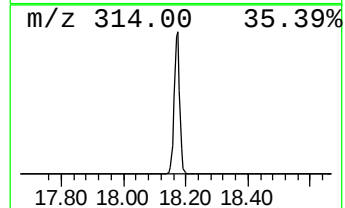
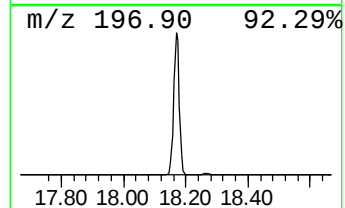
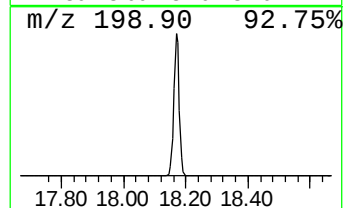
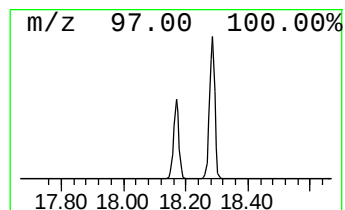
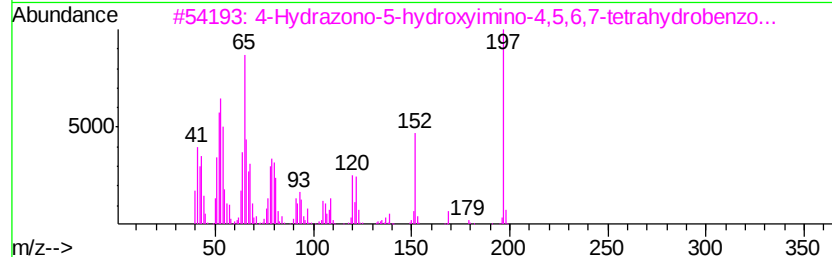
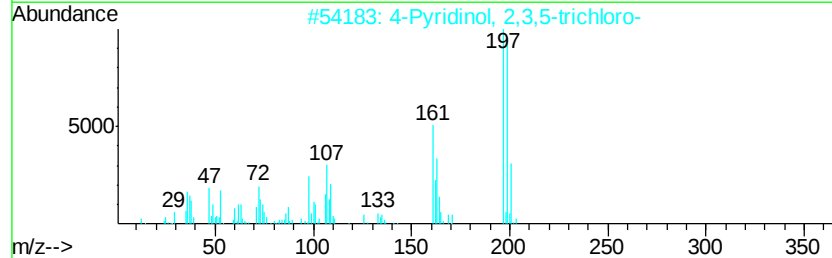
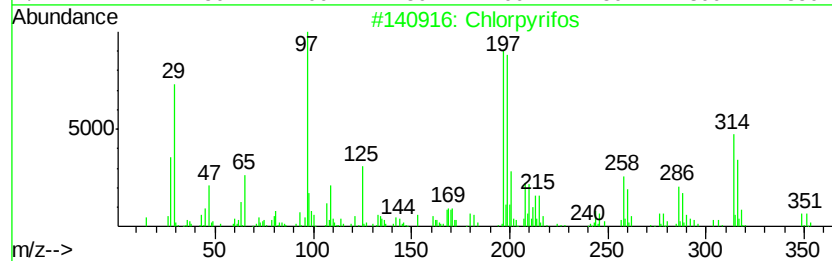
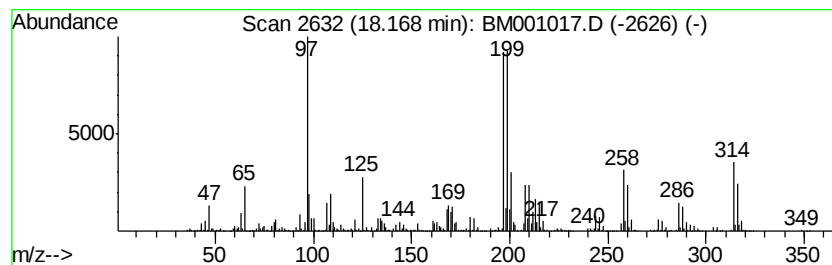
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Chlorpyrifos Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	10.05 ng/ul	3965510	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chlorpyrifos	349	C9H11Cl3N03PS	002921-88-2	91
2		4-Pyridinol, 2,3,5-trichloro-	197	C5H2Cl3NO	001970-40-7	27
3		4-Hydrazono-5-hydroxyimino-4,5,6...	197	C6H7N5O3	303194-86-7	25
4		2,4-Dichloroquinoline	197	C9H5Cl2N	000703-61-7	14
5		Formamide, N,N-diphenyl-	197	C13H11NO	000607-00-1	11



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
Data File : BM001017.D
Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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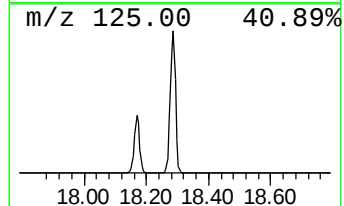
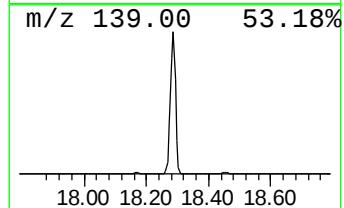
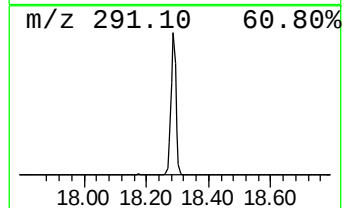
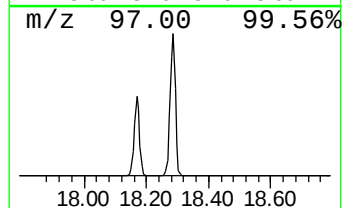
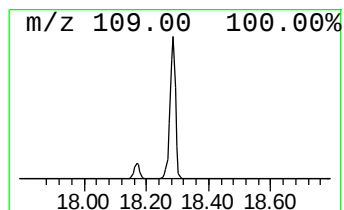
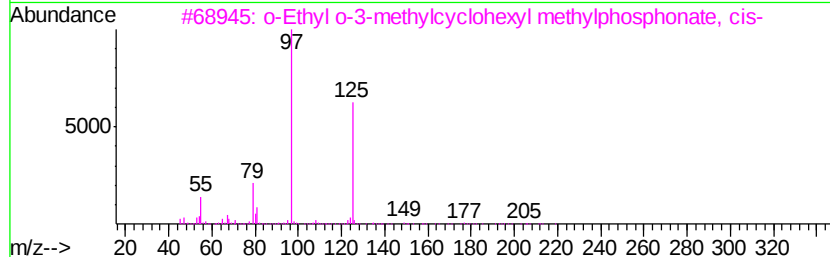
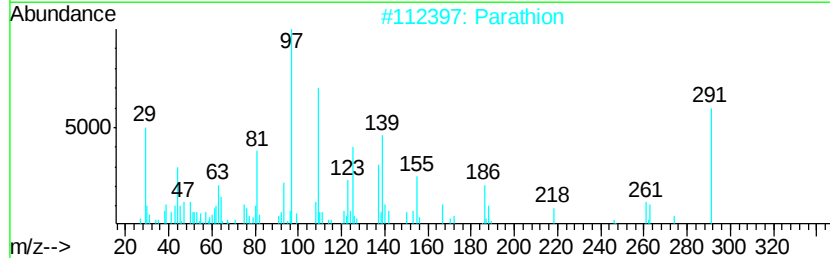
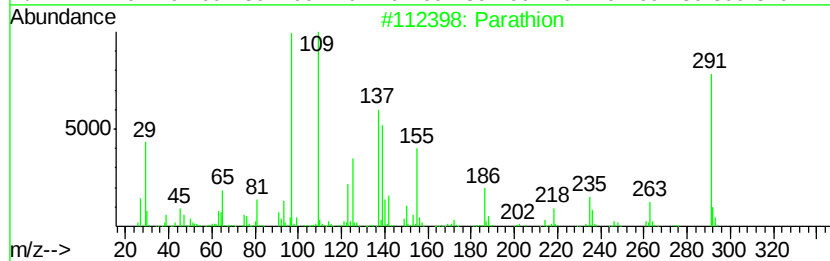
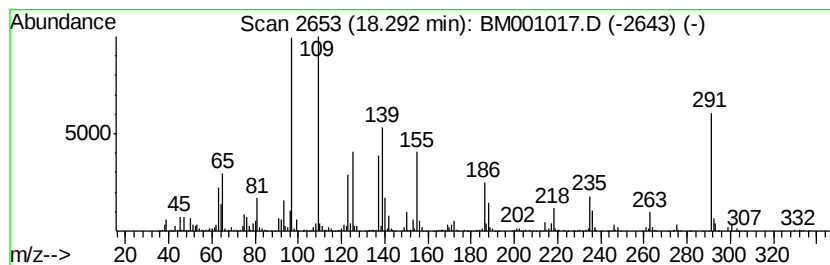
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Parathion Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.29	23.53 ng/ul	9284720	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Parathion	291	C10H14N05PS	000056-38-2	91
2		Parathion	291	C10H14N05PS	000056-38-2	90
3		o-Ethyl o-3-methylcyclohexyl met...	220	C10H2103P	1000275-28-1	22
4		o-Ethyl o-3-methylcyclohexyl met...	220	C10H2103P	1000275-28-2	12
5		o-Ethyl o-4-methylcyclohexyl met...	220	C10H2103P	1000275-28-4	12



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
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Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
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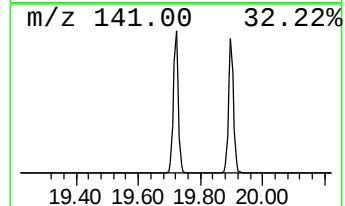
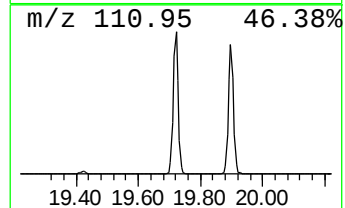
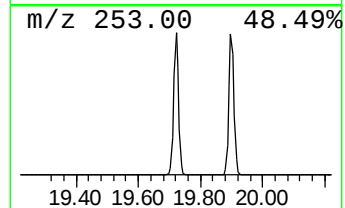
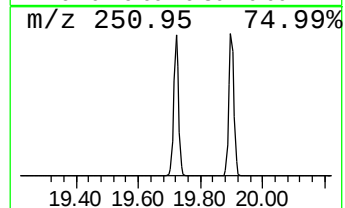
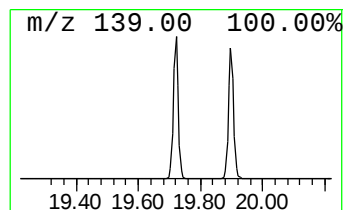
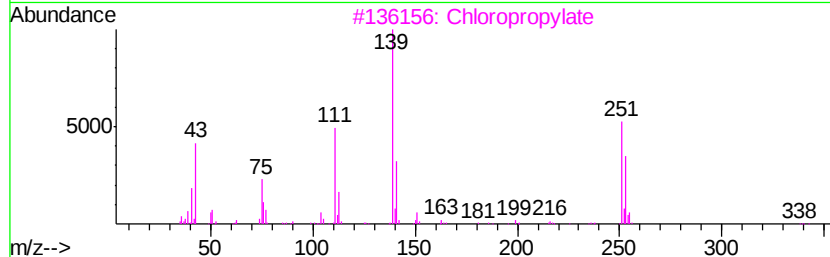
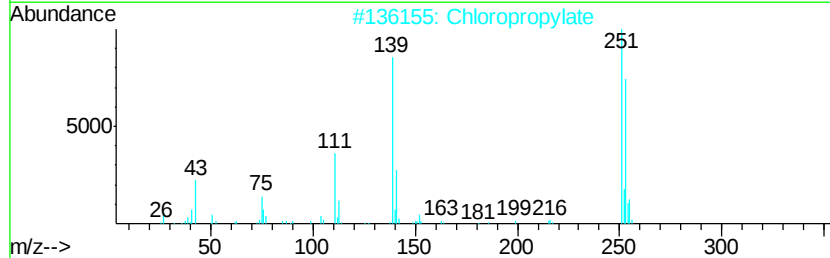
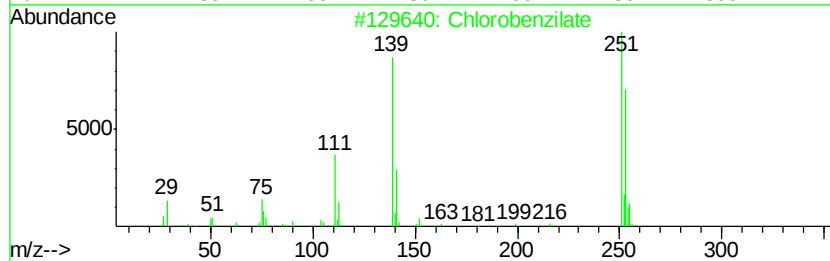
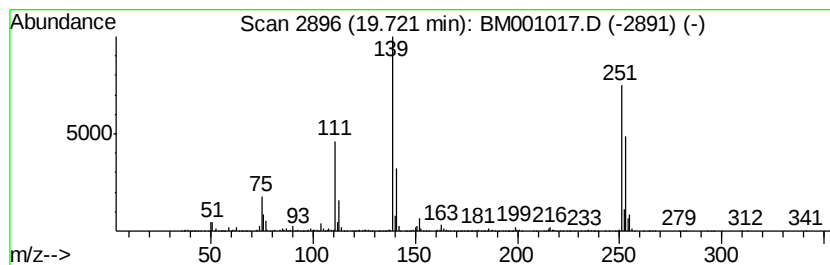
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM041515.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Chlorobenzilate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.72	28.60 ng/ul	5099730	Chrysene-d12	21.19
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Chlorobenzilate	324 C16H14Cl2O3	000510-15-6	94
2	Chloropropylate	338 C17H16Cl2O3	005836-10-2	91
3	Chloropropylate	338 C17H16Cl2O3	005836-10-2	91
4	Chlorobenzilate	324 C16H14Cl2O3	000510-15-6	91
5	4,4'-Dichloro-.alpha.-(trichloro...	368 C14H9Cl5O	000115-32-2	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
Data File : BM001017.D
Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
X1N61

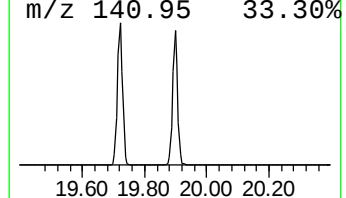
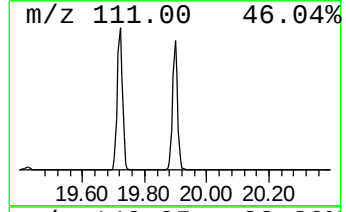
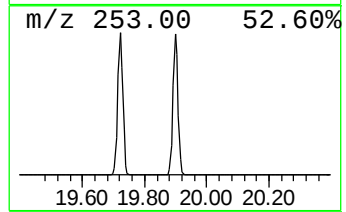
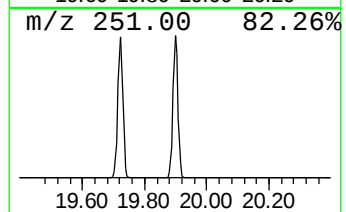
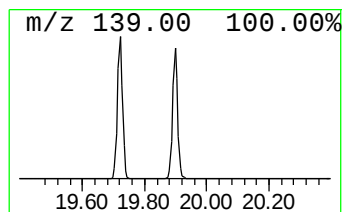
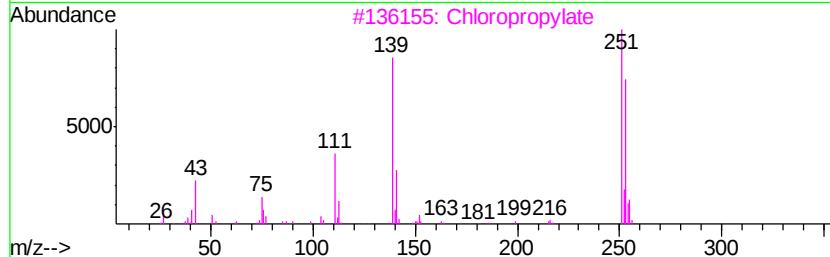
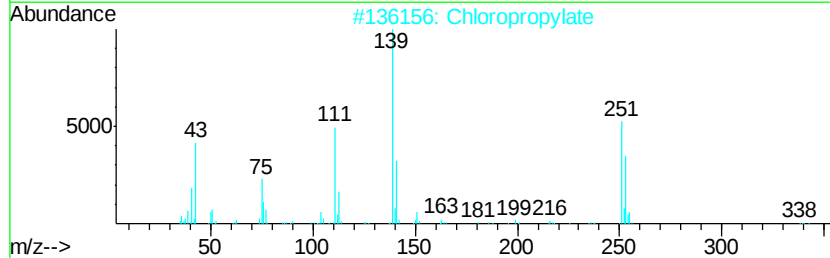
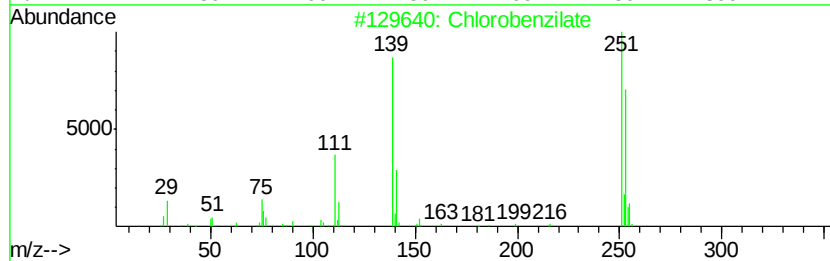
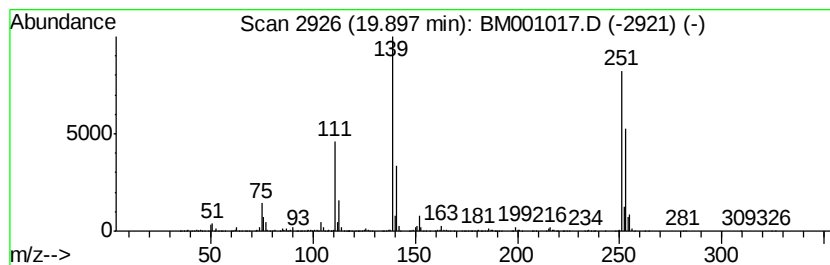
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4,4'-Dichloro-.alpha.-(tric... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.90	27.01 ng/ul	4816370	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chlorobenzilate	324	C16H14Cl2O3	000510-15-6	91
2		Chloropropylate	338	C17H16Cl2O3	005836-10-2	91
3		Chloropropylate	338	C17H16Cl2O3	005836-10-2	91
4		4,4'-Dichloro-.alpha.-(trichloro...	368	C14H9Cl5O	000115-32-2	91
5		Chlorobenzilate	324	C16H14Cl2O3	000510-15-6	86



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041615\
Data File : BM001017.D
Acq On : 16 Apr 2015 12:08
Operator : TP/IZ
Sample : G1749-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM041515.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
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(DEL) Alkane: Str...	7.80	3.2	ng/ul	304780	1	7.58	1902740	20.0
unknown-01	8.86	21.0	ng/ul	1997130	1	7.58	1902740	20.0
2-Hydroxy-3,5,6-t...	13.88	2.1	ng/ul	319786	3	14.24	3043950	20.0
Benzene, pentachl...	14.57	27.6	ng/ul	4207920	3	14.24	3043950	20.0
Chlorpyrifos	18.17	10.1	ng/ul	3965510	4	16.99	7891030	20.0
Parathion	18.29	23.5	ng/ul	9284720	4	16.99	7891030	20.0
Chlorobenzilate	19.72	28.6	ng/ul	5099730	5	21.19	3566120	20.0
4,4'-Dichloro-.al...	19.90	27.0	ng/ul	4816370	5	21.19	3566120	20.0