

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
 Data File : BF083019.D
 Acq On : 18 Nov 2015 20:29
 Operator : UM/IZ
 Sample : G4453-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CWC-MISC-1-20151116

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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_F\METHODS\8270-BF110715.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.201	181	185	189	rBV2	25930	49663	1.08%	0.087%
2	4.899	242	246	248	rBV	1546081	2617816	57.12%	4.592%
3	5.333	281	284	287	rBV	3444854	4233150	92.37%	7.425%
4	6.384	373	376	378	rBV	4086162	4582863	100.00%	8.039%
5	6.499	383	386	389	rBV	3966524	3904549	85.20%	6.849%
6	6.727	404	406	408	rBV	1169055	940145	20.51%	1.649%
7	6.887	417	420	422	rBV	3553347	3210860	70.06%	5.632%
8	7.299	452	456	458	rBV	2771599	3210380	70.05%	5.631%
9	8.019	516	519	521	rBV	1140888	1232431	26.89%	2.162%
10	9.093	610	613	616	rBV	4059776	4065621	88.71%	7.132%
11	9.493	645	648	650	rBV	972397	881812	19.24%	1.547%
12	9.768	669	672	674	rBV	1593192	1478548	32.26%	2.594%
13	10.213	710	711	713	rVB	73278	56721	1.24%	0.099%
14	10.419	727	729	732	rBV	250797	235660	5.14%	0.413%
15	10.556	738	741	743	rBV	2970387	3090525	67.44%	5.421%
16	10.682	751	752	756	rVB	67062	95560	2.09%	0.168%
17	10.751	756	758	762	rBV2	109020	102060	2.23%	0.179%
18	10.819	762	764	766	rVB	117685	103687	2.26%	0.182%
19	10.979	776	778	780	rBV	142093	158347	3.46%	0.278%
20	11.139	790	792	793	rVB	65640	75639	1.65%	0.133%
21	11.185	793	796	799	rVV	1193938	1513290	33.02%	2.654%
22	11.242	799	801	805	rVV2	1820798	2137594	46.64%	3.750%
23	11.345	808	810	812	rVB	646690	617939	13.48%	1.084%
24	11.505	818	824	827	rBV3	282484	606088	13.23%	1.063%
25	11.596	829	832	834	rVB	1531922	1918373	41.86%	3.365%
26	11.665	835	838	840	rVB	405593	411493	8.98%	0.722%
27	11.733	842	844	846	rBV2	552393	514548	11.23%	0.903%
28	11.768	846	847	850	rVB	230807	205154	4.48%	0.360%
29	11.825	850	852	853	rBV	80650	89292	1.95%	0.157%
30	11.871	853	856	857	rBV	508233	636449	13.89%	1.116%
31	11.905	857	859	860	rBV	284383	284938	6.22%	0.500%
32	12.031	868	870	874	rBV2	665659	1147688	25.04%	2.013%
33	12.088	874	875	877	rVV	263772	371599	8.11%	0.652%
34	12.134	877	879	881	rVB	364591	453708	9.90%	0.796%

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35	12.191	881	884	886	rBV2	115085	162971	3.56%	0.286%
36	12.294	890	893	894	rBV	74820	71025	1.55%	0.125%
37	12.328	894	896	897	rVV	385921	331625	7.24%	0.582%
38	12.385	897	901	903	rVB2	788924	1320120	28.81%	2.316%
39	12.454	903	907	909	rBV	493356	528131	11.52%	0.926%
40	12.511	910	912	914	rBV	320171	484165	10.56%	0.849%
41	12.556	914	916	918	rVV2	126912	161893	3.53%	0.284%
42	12.602	918	920	922	rVB2	51122	69842	1.52%	0.123%
43	12.671	924	926	929	rBV	369988	478614	10.44%	0.840%
44	12.728	929	931	932	rVB2	50669	62504	1.36%	0.110%
45	12.774	933	935	938	rBV2	73649	106433	2.32%	0.187%
46	12.831	938	940	943	rVV	3500120	3650918	79.66%	6.404%
47	12.876	943	944	948	rVV3	45859	73739	1.61%	0.129%
48	13.002	952	955	958	rVB	42829	92374	2.02%	0.162%
49	13.048	958	959	960	rBV	104347	88801	1.94%	0.156%
50	13.105	963	964	970	rVB3	58724	87280	1.90%	0.153%
51	13.196	970	972	974	rBV2	40254	55592	1.21%	0.098%
52	13.254	974	977	978	rBV2	71834	80184	1.75%	0.141%
53	13.391	985	989	990	rBV2	26104	64216	1.40%	0.113%
54	13.505	997	999	1002	rBV	38112	50002	1.09%	0.088%
55	13.619	1006	1009	1011	rBV	50127	83482	1.82%	0.146%
56	13.722	1016	1018	1019	rBV	29576	47167	1.03%	0.083%
57	13.745	1019	1020	1023	rVB2	63976	95539	2.08%	0.168%
58	13.871	1028	1031	1037	rVB2	1282120	1522366	33.22%	2.670%
59	14.877	1116	1119	1124	rVB	106019	224487	4.90%	0.394%
60	15.140	1140	1142	1145	rBV	53129	77956	1.70%	0.137%
61	15.197	1145	1147	1150	rBV	93595	120480	2.63%	0.211%
62	15.254	1150	1152	1155	rBV	535479	789346	17.22%	1.385%
63	16.557	1264	1266	1270	rVB2	42135	90404	1.97%	0.159%
64	16.945	1297	1300	1304	rBV	27660	62746	1.37%	0.110%
65	18.146	1401	1405	1411	rVB4	24185	77744	1.70%	0.136%
66	20.089	1567	1575	1586	rVB2	110693	496634	10.84%	0.871%
67	21.174	1664	1670	1671	rBV4	20577	65977	1.44%	0.116%

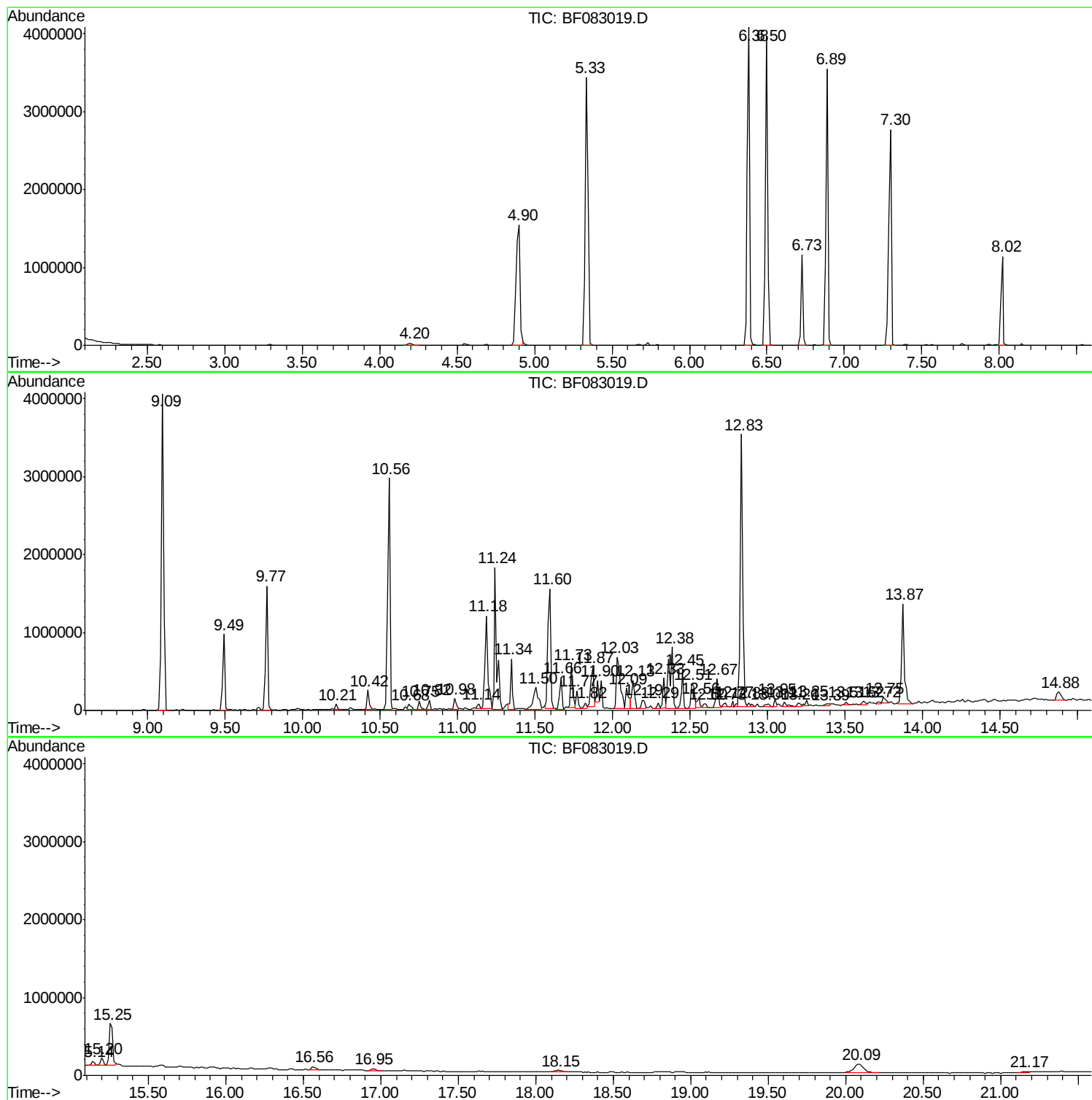
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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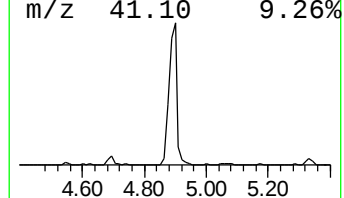
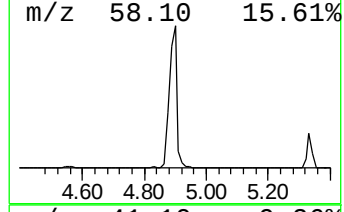
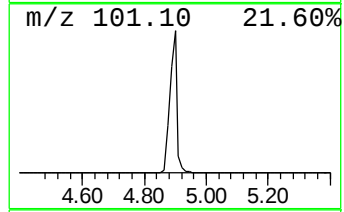
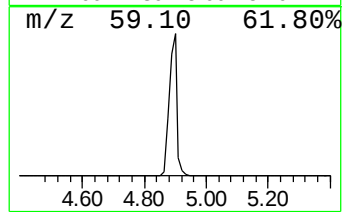
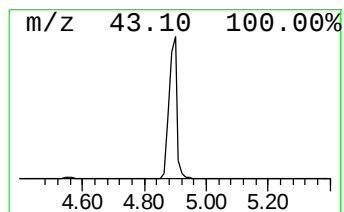
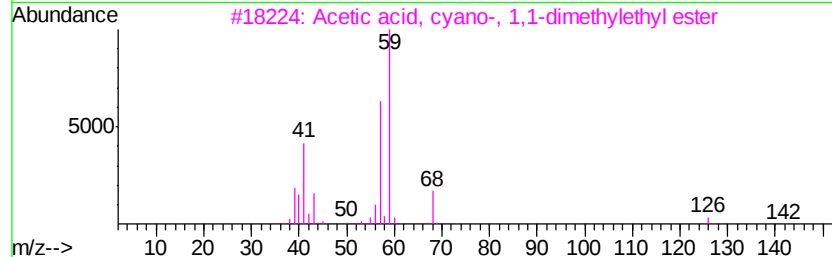
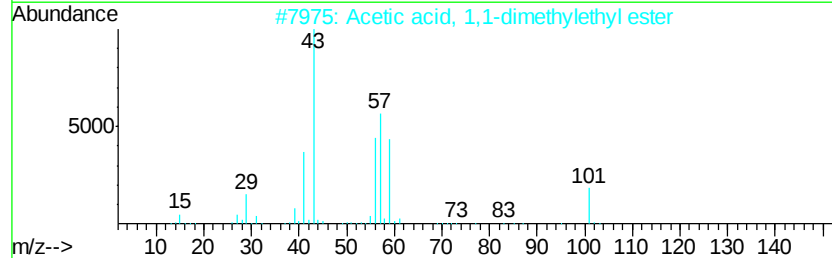
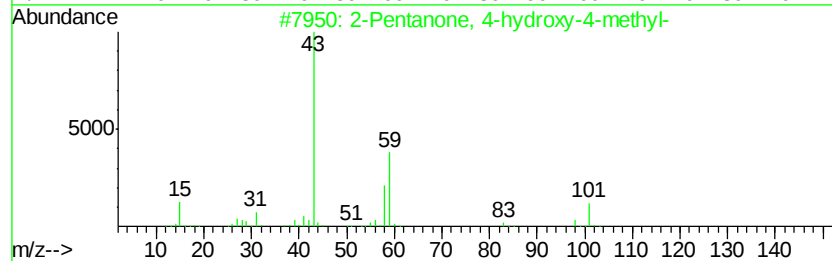
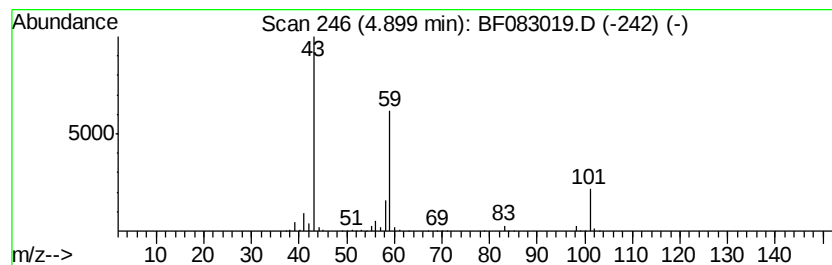
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	55.69 ng	2617820	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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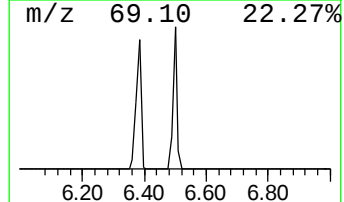
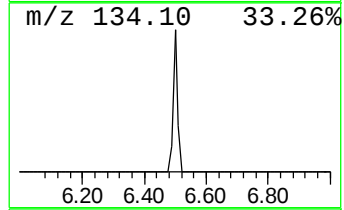
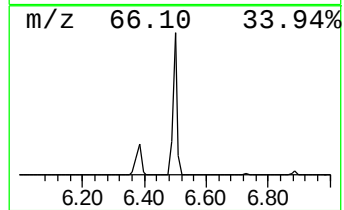
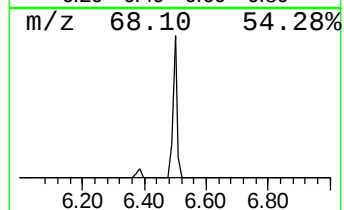
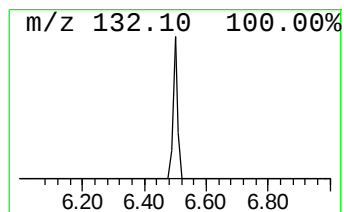
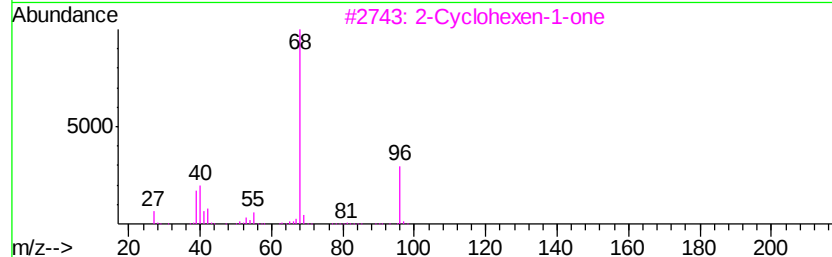
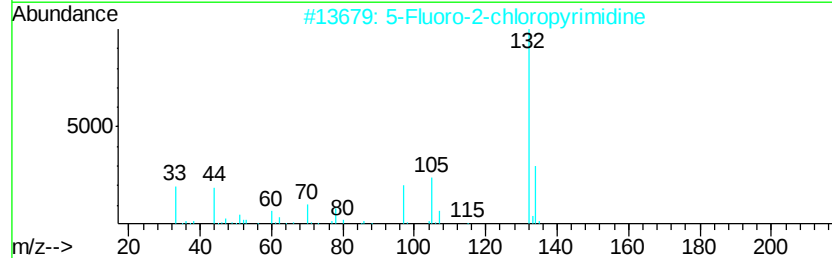
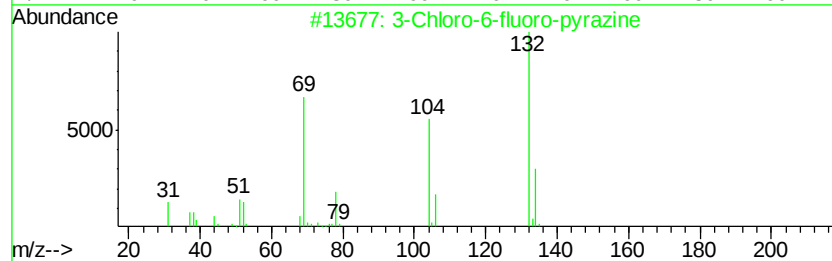
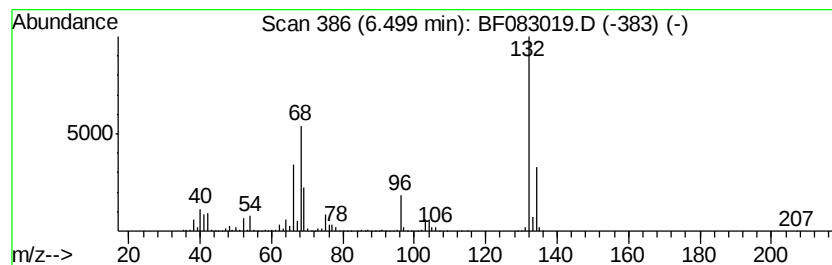
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	83.06 ng	3904550	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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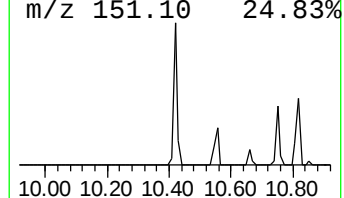
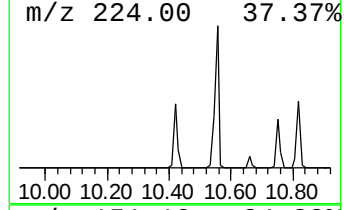
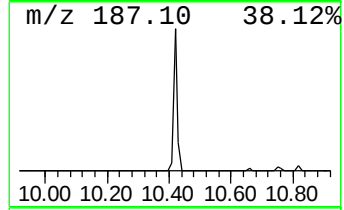
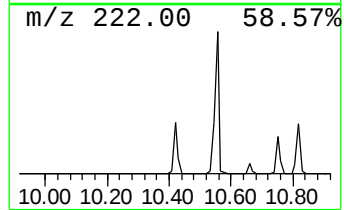
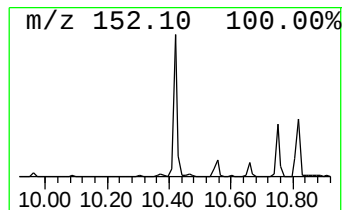
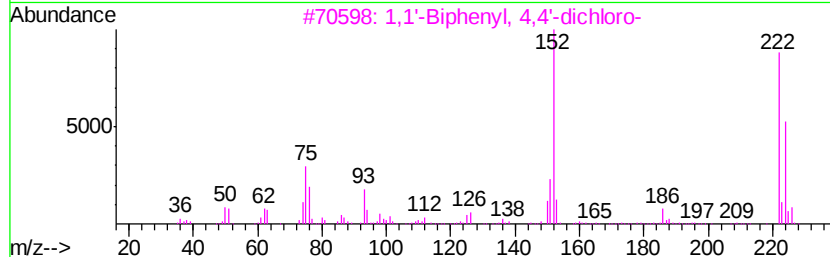
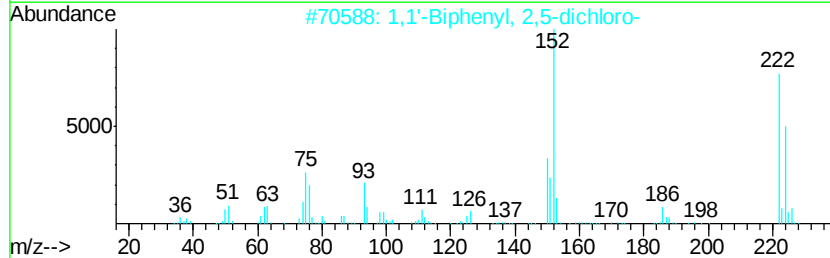
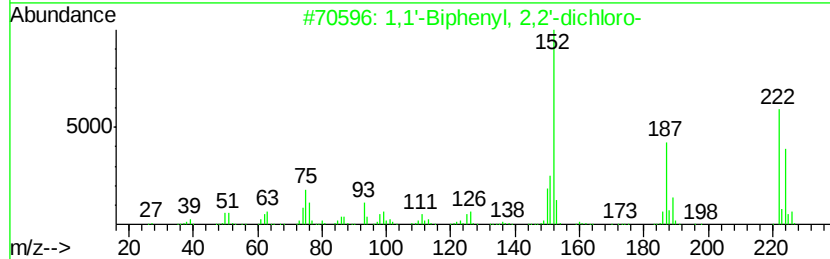
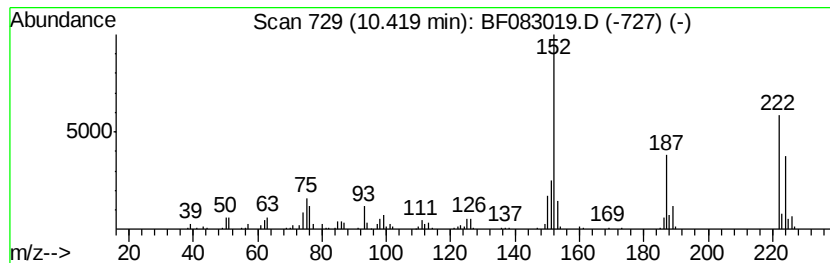
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.42	3.19 ng	235660	Acenaphthene-d10	9.77	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2	1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	95
3	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	95
4	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	95
5	1,1'-Biphenyl, 2,6-dichloro-	222	C12H8Cl2	033146-45-1	89



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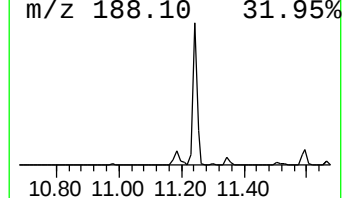
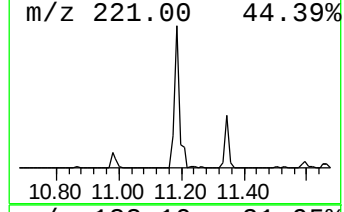
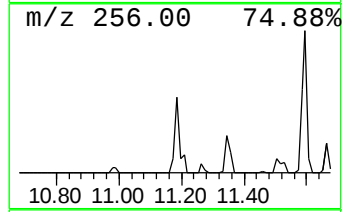
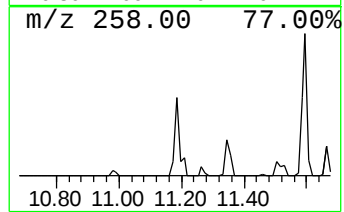
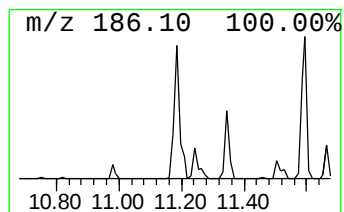
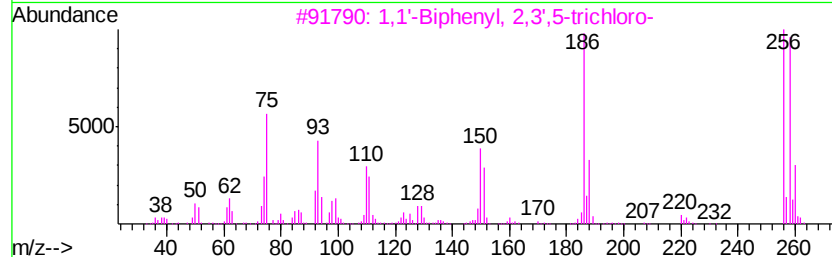
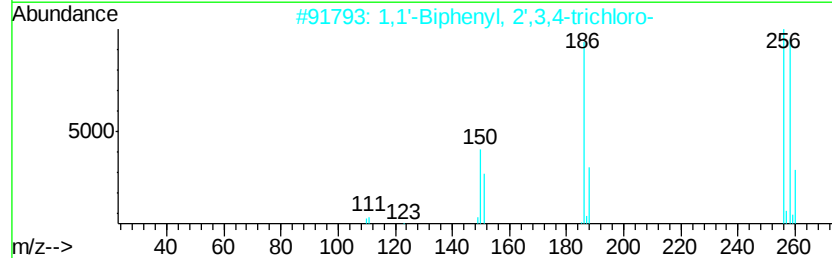
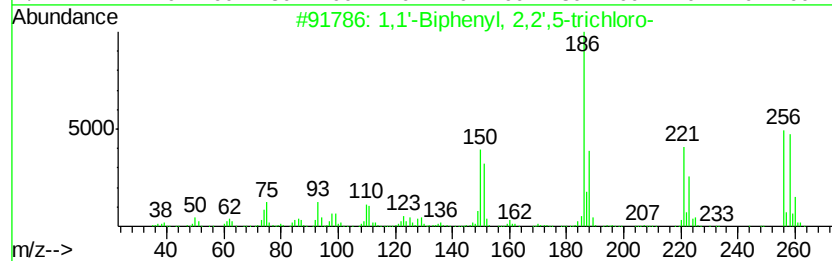
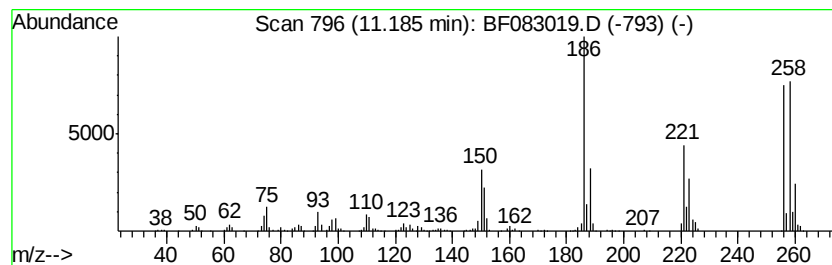
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Peak Number 5 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.18	14.16 ng	1513290	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	96
3	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	93
4	1,1'-Biphenyl, 3,4,4'-Trichloro-	256	C12H7Cl3	038444-90-5	92
5	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	90



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Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

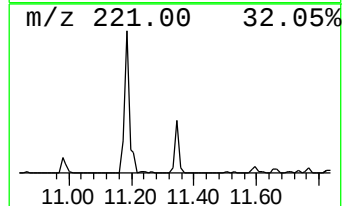
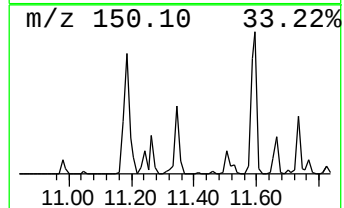
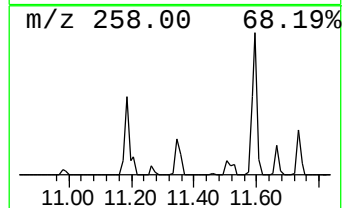
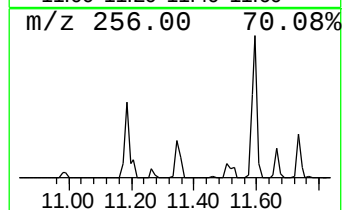
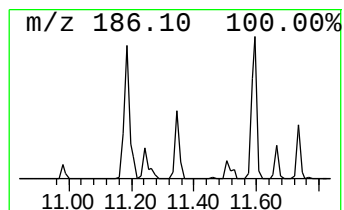
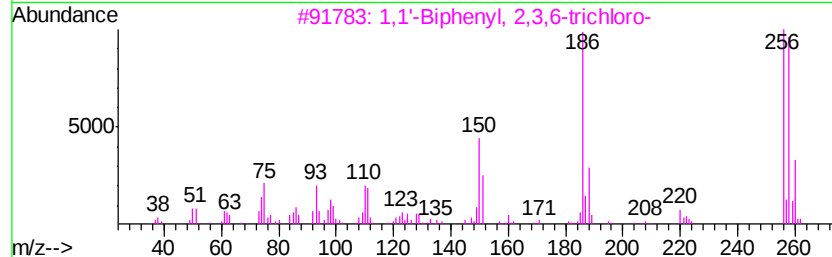
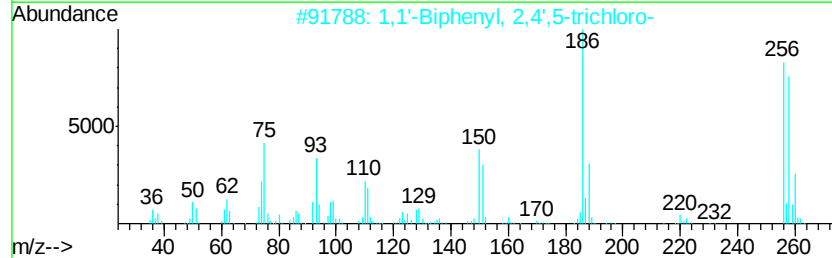
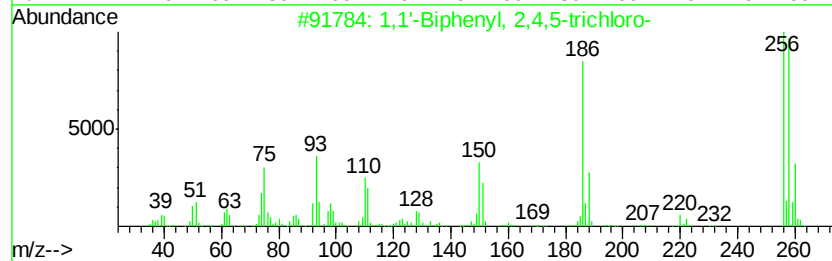
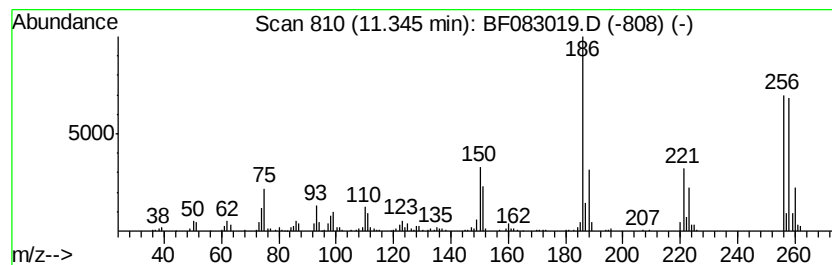
Instrument :
BNA_F
ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,4,5-trichl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.34	5.78 ng	617939	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98
2	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
3	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	96
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	96
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083019.D
Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

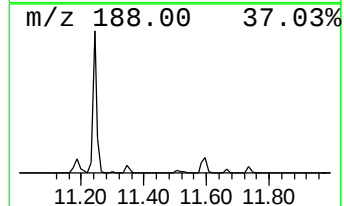
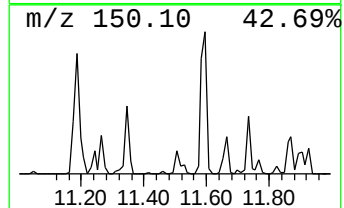
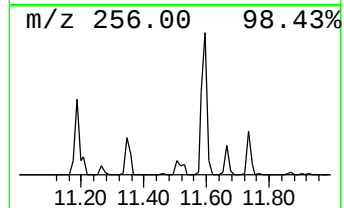
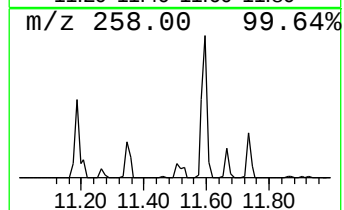
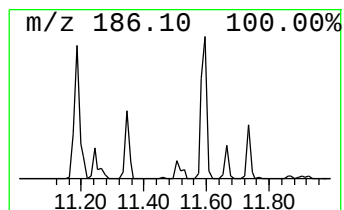
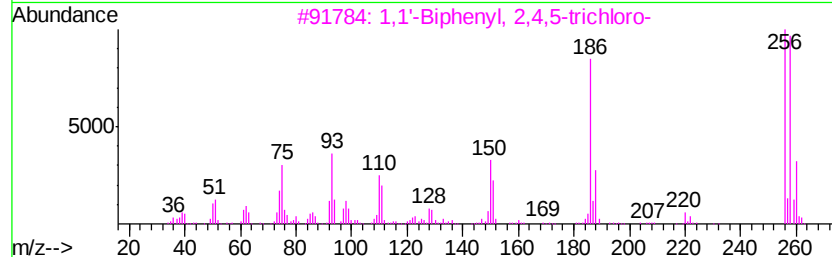
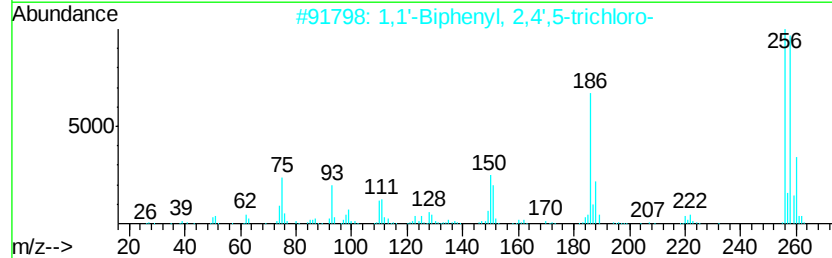
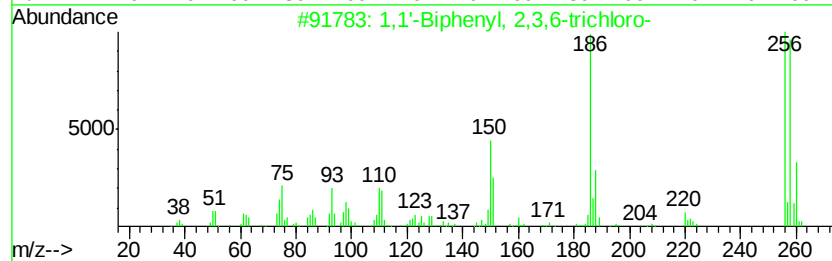
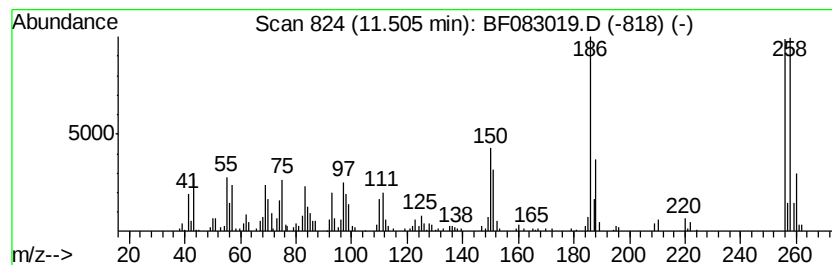
Instrument :
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ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.50	5.67 ng	606088	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99
2	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
3	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98
4	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	98
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	97



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

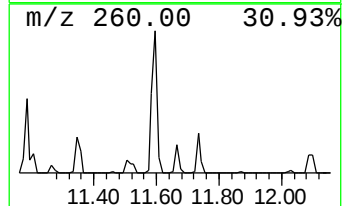
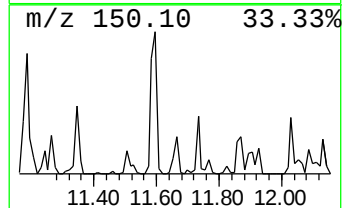
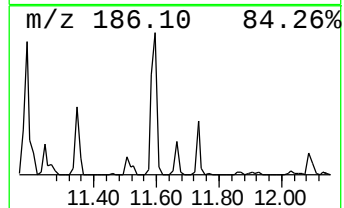
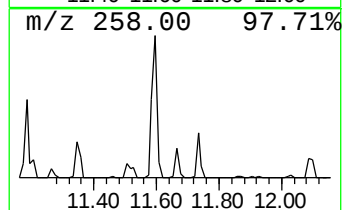
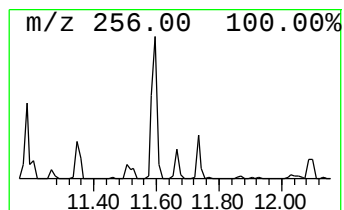
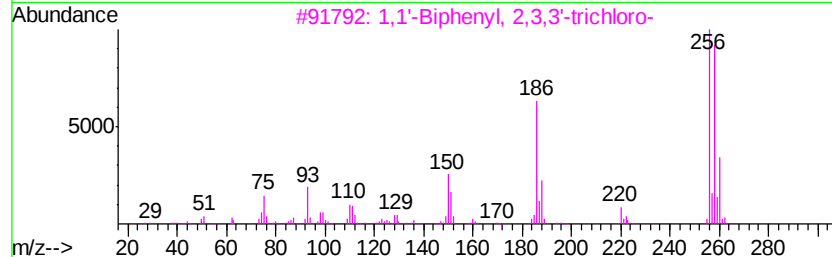
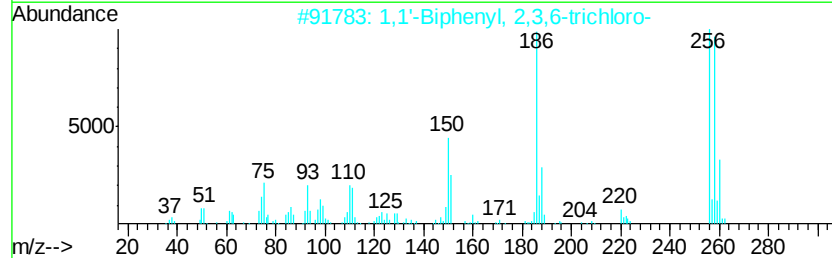
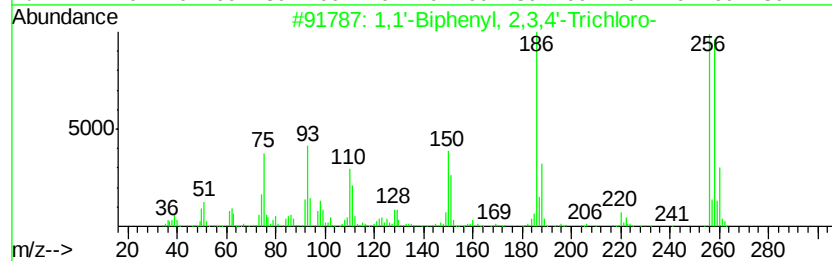
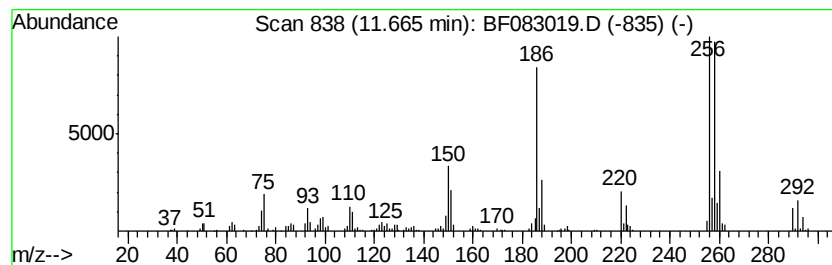
Instrument :
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ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1,1'-Biphenyl, 2,3,4'-Trich... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.66	3.85 ng	411493	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	98
2	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	98
3	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	98
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	98
5	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083019.D
Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

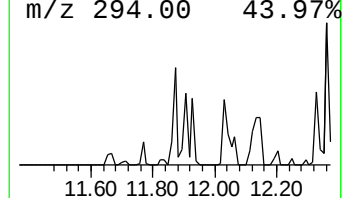
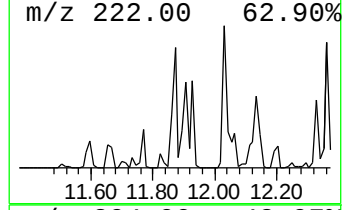
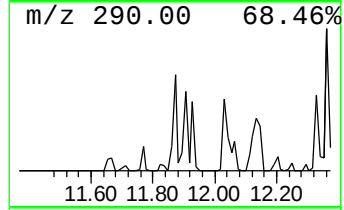
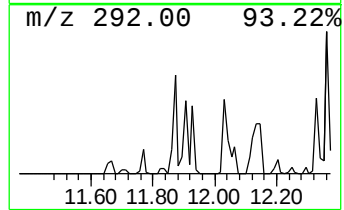
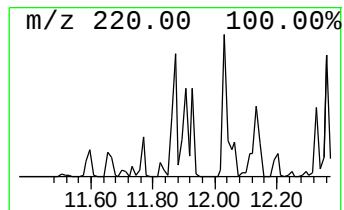
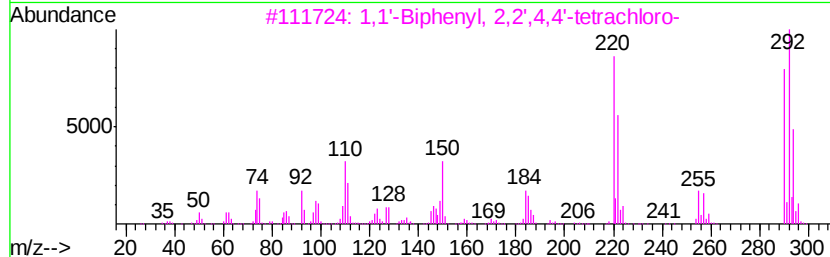
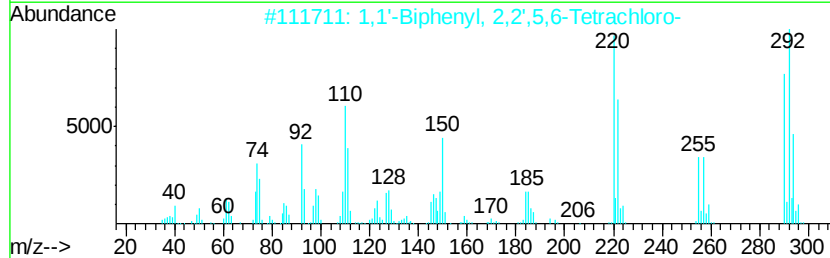
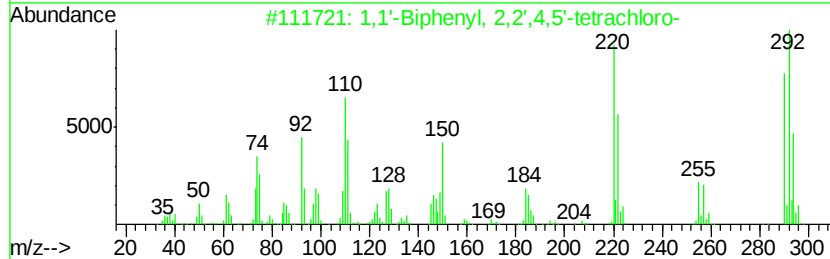
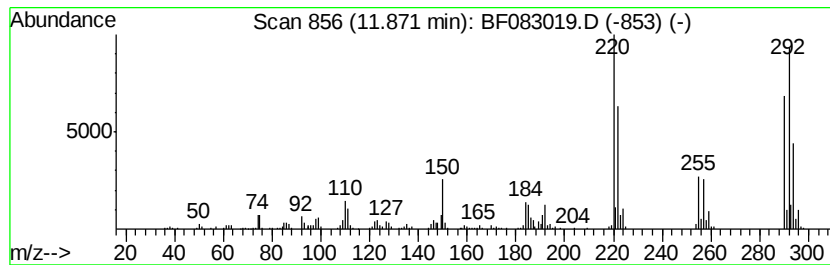
Instrument :
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ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.87	5.95 ng	636449	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
2	1,1'-Biphenyl, 2,2',5,6-Tetrachl...	290	C12H6Cl4	041464-41-9	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	96
5	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083019.D
Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

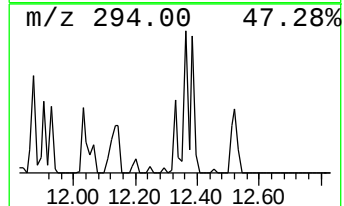
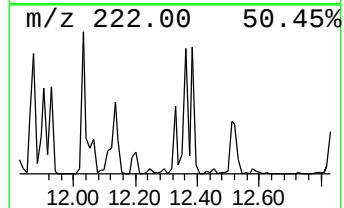
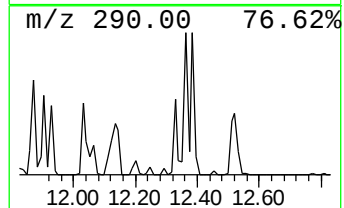
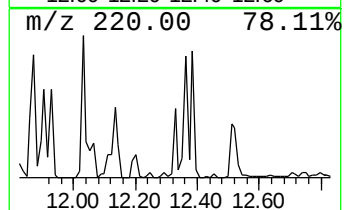
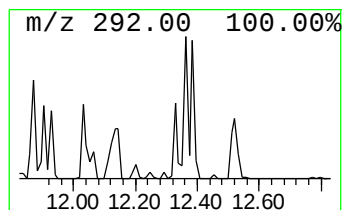
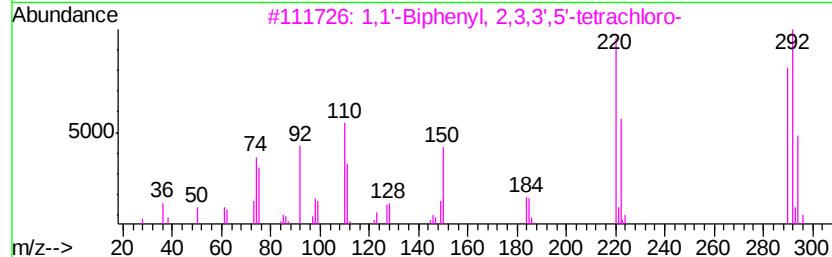
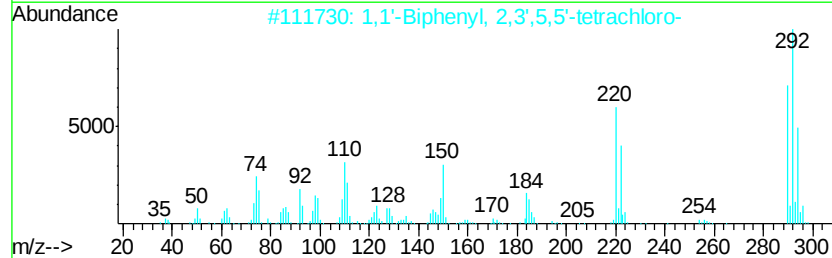
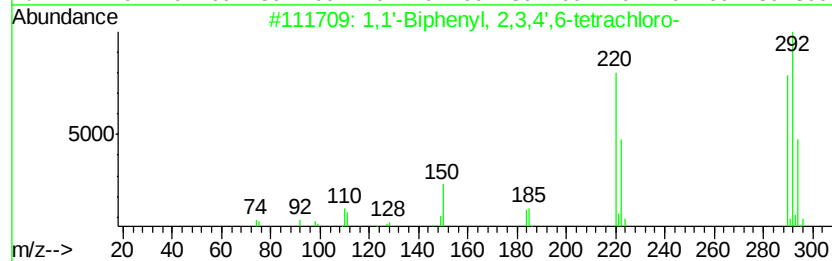
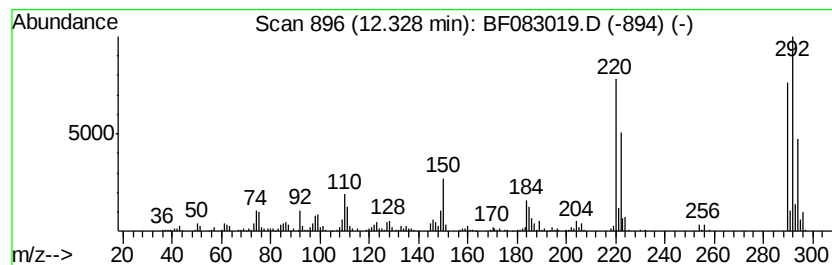
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CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1,1'-Biphenyl, 2,3,4',6-tet... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.33	3.10 ng	331625	Phenanthrene-d10	11.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99
2	1,1'-Biphenyl,	2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
3	1,1'-Biphenyl,	2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
4	1,1'-Biphenyl,	2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	1,1'-Biphenyl,	2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083019.D
Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

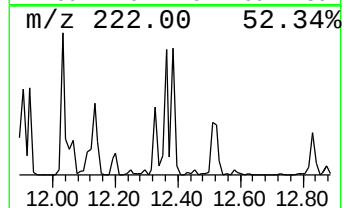
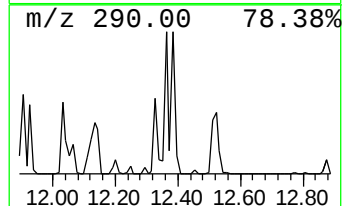
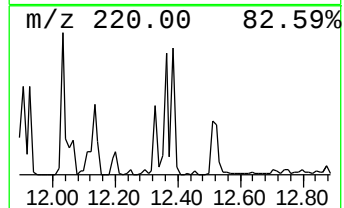
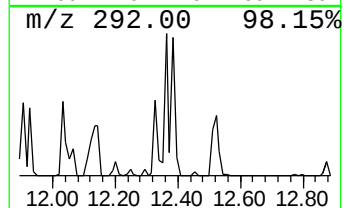
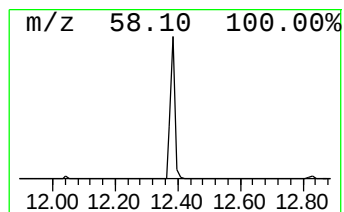
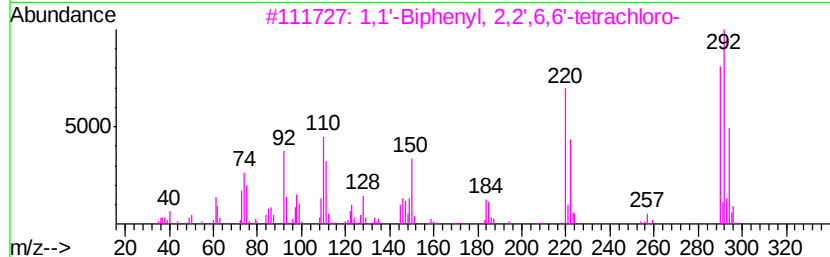
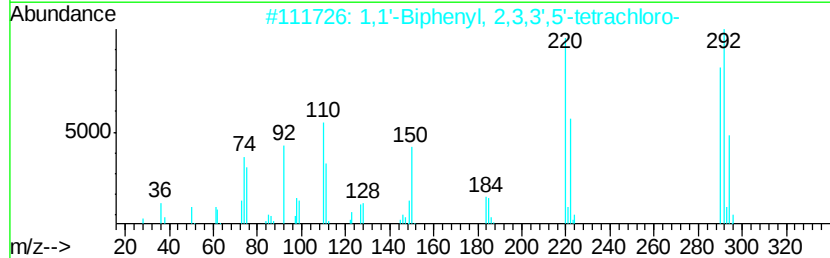
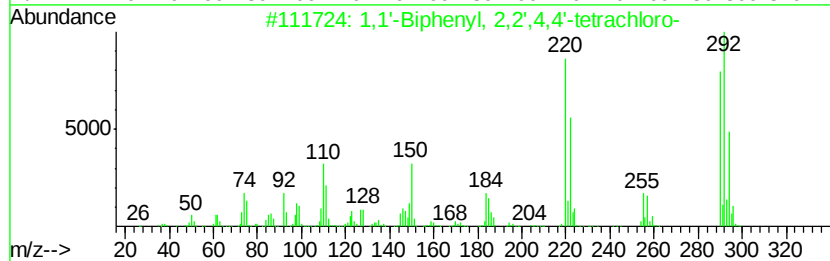
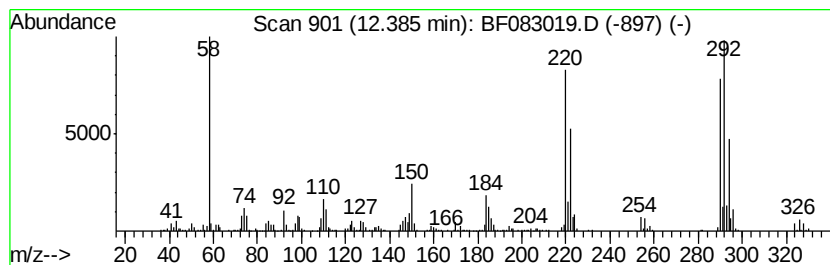
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.39	12.35 ng	1320120	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
4	1,1'-Biphenyl, 2,3',4',5'-tetrach...	290	C12H6Cl4	032598-11-1	98
5	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
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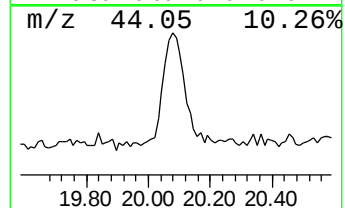
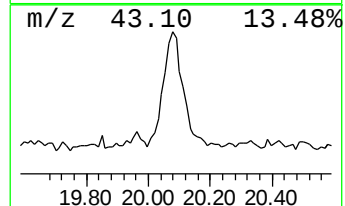
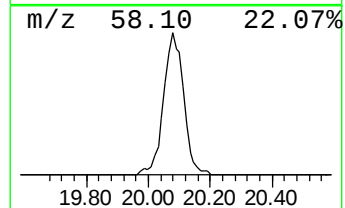
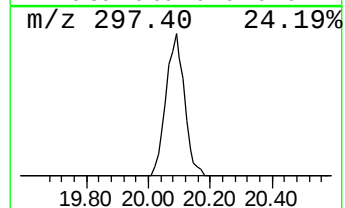
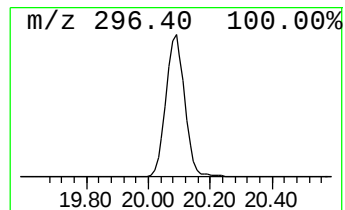
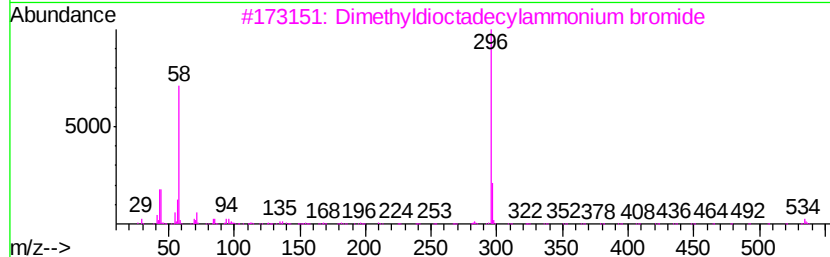
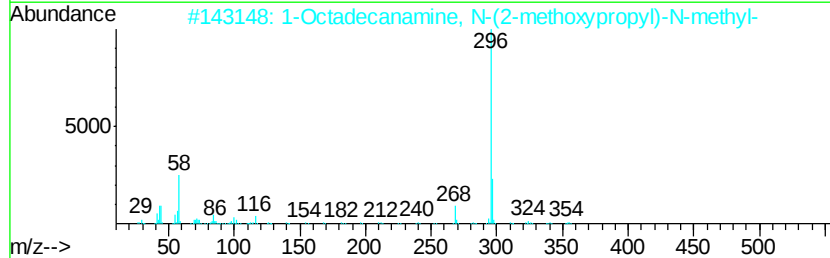
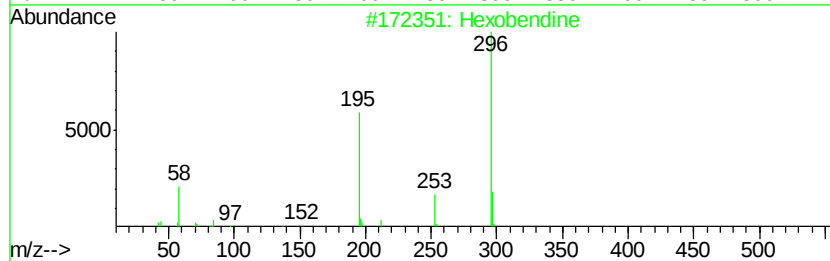
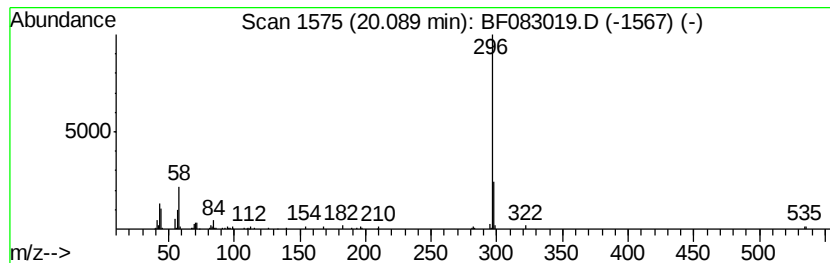
Instrument :
BNA_F
ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Hexobendine Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.09	12.58 ng	496634	Perylene-d12	15.27
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexobendine	592	C30H44N2O10	000054-03-5 78
2	1-Octadecanamine, N-(2-methoxypr...	355	C23H49NO	1000196-18-5 78
3	Dimethyldioctadecylammonium bromide	630	C38H80BrN	003700-67-2 64
4	Benzeneethanamine, .beta.-hydrox...	403	C27H49NO	1000197-20-9 59
5	2-(2-Quinolyl)(1,2,4)triazolo(1,...	296	C19H12N4	1000241-15-8 45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083019.D
Acq On : 18 Nov 2015 20:29
Operator : UM/IZ
Sample : G4453-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CWC-MISC-1-20151116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
2-Pentanone, 4-hy...	4.90	55.7	ng	2617820	1	6.73	940145	20.0
unknown6.50	6.50	83.1	ng	3904550	1	6.73	940145	20.0
1,1'-Biphenyl, 2,...	10.42	3.2	ng	235660	3	9.77	1478550	20.0
1,1'-Biphenyl, 2,...	11.18	14.2	ng	1513290	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	11.34	5.8	ng	617939	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	11.50	5.7	ng	606088	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	11.66	3.9	ng	411493	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	11.87	6.0	ng	636449	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	12.33	3.1	ng	331625	4	11.24	2137590	20.0
1,1'-Biphenyl, 2,...	12.39	12.4	ng	1320120	4	11.24	2137590	20.0
Hexobendine	20.09	12.6	ng	496634	6	15.27	789346	20.0