

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
 Data File : BF082445.D
 Acq On : 22 Oct 2015 11:57
 Operator : UM/IZ
 Sample : G4111-03
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CLEARWELL

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-BF101015.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	6.537	381	385	387	rBV2	1436204	1721642	7.21%	0.944%
2	6.926	411	419	420	rBV	8326155	23867932	100.00%	13.082%
3	7.349	453	456	457	rBV	879717	1245443	5.22%	0.683%
4	7.600	473	478	481	rBV	618060	855507	3.58%	0.469%
5	7.886	495	503	505	rVV2	390783	1278158	5.36%	0.701%
6	8.058	516	518	519	rVV	880117	756144	3.17%	0.414%
7	9.143	610	613	615	rBV	2628460	2635187	11.04%	1.444%
8	9.818	669	672	679	rVB2	858946	1096923	4.60%	0.601%
9	10.275	710	712	716	rVV4	339300	788836	3.31%	0.432%
10	10.389	719	722	728	rVB	694253	1616006	6.77%	0.886%
11	10.606	739	741	746	rBV2	796894	891904	3.74%	0.489%
12	10.732	750	752	756	rBV2	1005855	1898396	7.95%	1.040%
13	10.801	756	758	760	rVB	968056	841521	3.53%	0.461%
14	10.892	764	766	770	rVB	1147992	1124859	4.71%	0.617%
15	10.972	770	773	775	rBV	832884	1096498	4.59%	0.601%
16	11.178	786	791	792	rBV2	1334208	2763464	11.58%	1.515%
17	11.304	799	802	804	rVB	1129765	1575239	6.60%	0.863%
18	11.338	804	805	810	rVB	561689	850853	3.56%	0.466%
19	11.418	810	812	816	rBV	1994129	3235339	13.56%	1.773%
20	11.486	816	818	822	rVB	1271960	1322078	5.54%	0.725%
21	11.589	825	827	829	rBV	1709304	2194445	9.19%	1.203%
22	11.681	833	835	839	rVB	1240050	1227108	5.14%	0.673%
23	11.818	845	847	849	rBV2	1912028	3091411	12.95%	1.694%
24	11.852	849	850	852	rVB	1415142	1074732	4.50%	0.589%
25	11.898	852	854	859	rVB	1686712	2367765	9.92%	1.298%
26	12.001	861	863	868	rVB	1822098	1802611	7.55%	0.988%
27	12.092	868	871	875	rBV	1627650	1744747	7.31%	0.956%
28	12.207	877	881	883	rBV	917812	1883742	7.89%	1.032%
29	12.241	883	884	886	rVB	882537	718854	3.01%	0.394%
30	12.287	886	888	892	rVB2	1121714	1712312	7.17%	0.938%
31	12.401	895	898	903	rVB	1063844	1975664	8.28%	1.083%
32	12.481	903	905	909	rBV	1695097	3439030	14.41%	1.885%
33	12.549	909	911	913	rBV	1275320	1286550	5.39%	0.705%
34	12.652	918	920	921	rBV	1248549	1420137	5.95%	0.778%

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35	12.744	926	928	931	rVV	995375	1173062	4.91%	0.643%
36	12.835	934	936	938	rBV	1912843	3051768	12.79%	1.673%
37	12.869	938	939	941	rVV	1727947	1614644	6.76%	0.885%
38	12.927	941	944	950	rVB2	1379005	3828114	16.04%	2.098%
39	13.029	950	953	958	rVB	1194032	2037320	8.54%	1.117%
40	13.121	958	961	964	rBV	1305578	1872579	7.85%	1.026%
41	13.201	964	968	970	rBV	918929	1931650	8.09%	1.059%
42	13.235	970	971	974	rVB	832345	851268	3.57%	0.467%
43	13.292	974	976	978	rBV	734534	999953	4.19%	0.548%
44	13.327	978	979	981	rVV	1310442	1311039	5.49%	0.719%
45	13.395	983	985	986	rVV	976028	1219598	5.11%	0.668%
46	13.464	986	991	992	rVV	4561396	9412226	39.43%	5.159%
47	13.544	996	998	1000	rVV	957356	1128980	4.73%	0.619%
48	13.647	1004	1007	1009	rVV2	977038	1541077	6.46%	0.845%
49	13.681	1009	1010	1013	rVB	1237579	1250106	5.24%	0.685%
50	13.738	1013	1015	1018	rVB	726434	1043858	4.37%	0.572%
51	13.807	1018	1021	1022	rBV	1799459	2540837	10.65%	1.393%
52	13.841	1022	1024	1027	rVV	1314988	1924929	8.06%	1.055%
53	13.898	1027	1029	1031	rVB	1014720	1446245	6.06%	0.793%
54	13.944	1031	1033	1036	rVB	1006176	1166147	4.89%	0.639%
55	14.001	1036	1038	1039	rBV	1249151	1536387	6.44%	0.842%
56	14.024	1039	1040	1044	rVB	615276	886686	3.71%	0.486%
57	14.092	1044	1046	1048	rBV	1354993	1643750	6.89%	0.901%
58	14.138	1048	1050	1052	rVV	731797	1403341	5.88%	0.769%
59	14.184	1052	1054	1057	rVB	604183	813701	3.41%	0.446%
60	14.298	1061	1064	1067	rVV	1672769	2280080	9.55%	1.250%
61	14.355	1067	1069	1072	rVB	539524	844853	3.54%	0.463%
62	14.447	1075	1077	1079	rVV	1368967	1796142	7.53%	0.984%
63	14.492	1079	1081	1083	rVV	1738406	2544513	10.66%	1.395%
64	14.607	1089	1091	1093	rVB	716267	1041715	4.36%	0.571%
65	14.653	1093	1095	1097	rBV	1552838	1764284	7.39%	0.967%
66	14.698	1097	1099	1101	rVB	617235	825363	3.46%	0.452%
67	14.744	1101	1103	1105	rBV	1528305	2440166	10.22%	1.337%
68	14.778	1105	1106	1109	rVV	1333039	1391888	5.83%	0.763%
69	14.835	1109	1111	1115	rVB	1235507	1599035	6.70%	0.876%
70	14.904	1115	1117	1119	rBV	1098661	1428386	5.98%	0.783%
71	14.950	1119	1121	1123	rBV	1018240	1251685	5.24%	0.686%

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72	15.041	1127	1129	1130	rBV	1183470	1536319	6.44%	0.842%
73	15.178	1138	1141	1143	rBV	494029	801534	3.36%	0.439%
74	15.247	1145	1147	1150	rVV	1448880	2181012	9.14%	1.195%
75	15.350	1153	1156	1157	rBV	699251	946986	3.97%	0.519%
76	15.384	1157	1159	1162	rVV	722880	1490468	6.24%	0.817%
77	15.430	1162	1163	1168	rVB	586670	1023071	4.29%	0.561%
78	15.624	1178	1180	1182	rVV	985966	1481336	6.21%	0.812%
79	15.704	1185	1187	1189	rVV	931223	1613066	6.76%	0.884%
80	15.750	1189	1191	1194	rVB	714836	1198090	5.02%	0.657%
81	15.818	1194	1197	1202	rVB	768869	1390850	5.83%	0.762%
82	15.910	1202	1205	1207	rBV	722137	1213555	5.08%	0.665%
83	15.955	1207	1209	1213	rVB	660758	1049347	4.40%	0.575%
84	16.081	1217	1220	1225	rBV	878320	2262462	9.48%	1.240%
85	16.253	1231	1235	1241	rBV	253994	711166	2.98%	0.390%
86	16.367	1242	1245	1248	rBV	789415	1679406	7.04%	0.920%
87	16.493	1253	1256	1258	rVV	452844	1139564	4.77%	0.625%
88	16.527	1258	1259	1262	rVV	456806	1046725	4.39%	0.574%
89	16.618	1265	1267	1274	rVV	403129	1296770	5.43%	0.711%
90	16.801	1280	1283	1288	rVV	281612	725257	3.04%	0.397%
91	16.904	1288	1292	1295	rVV	616140	1516741	6.35%	0.831%
92	16.984	1295	1299	1304	rVV2	1096591	3667003	15.36%	2.010%
93	17.064	1304	1306	1312	rVV	429451	1064934	4.46%	0.584%
94	17.167	1312	1315	1325	rVB	347574	1101729	4.62%	0.604%
95	17.338	1325	1330	1332	rBV	312810	871785	3.65%	0.478%
96	17.384	1332	1334	1340	rVB	272422	723022	3.03%	0.396%
97	17.578	1346	1351	1357	rBV	378750	1428230	5.98%	0.783%
98	18.024	1385	1390	1400	rBV	323387	1267119	5.31%	0.694%
99	18.847	1456	1462	1466	rBV	195152	772501	3.24%	0.423%
100	18.950	1466	1471	1478	rVV2	195885	1016592	4.26%	0.557%

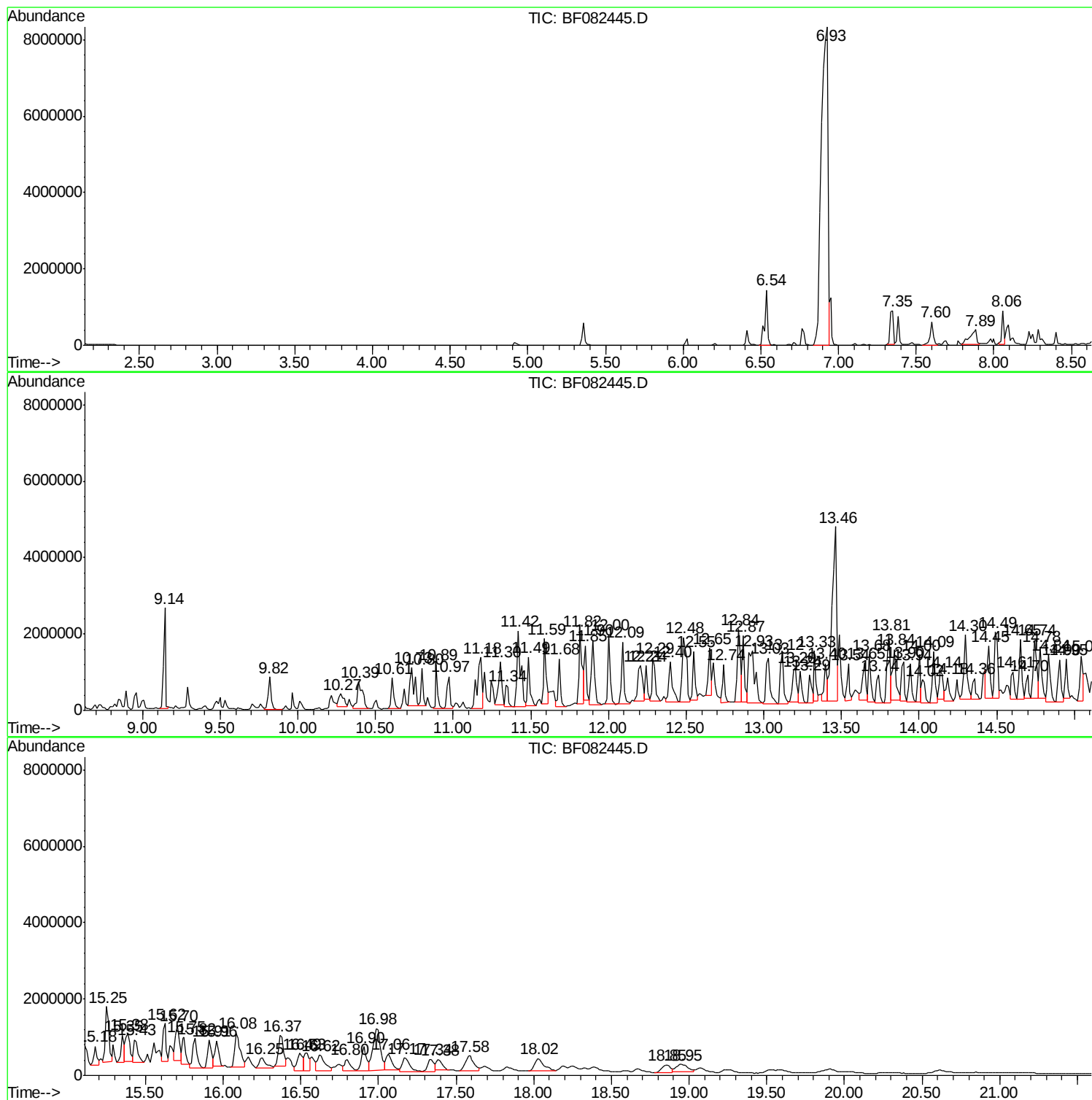
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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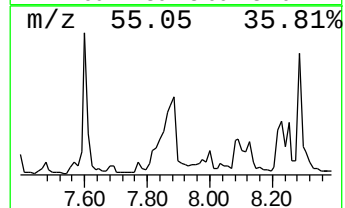
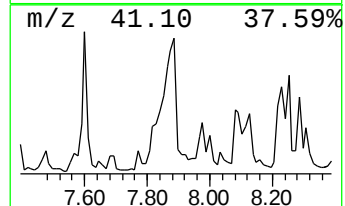
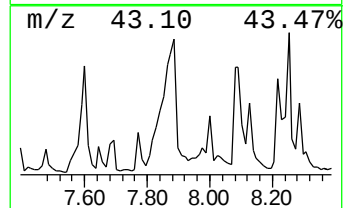
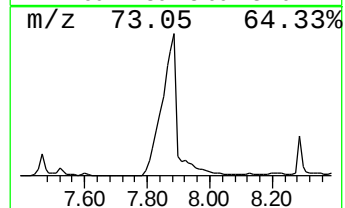
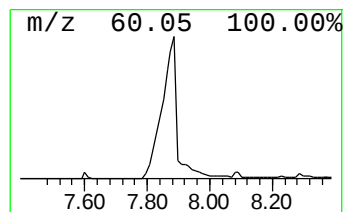
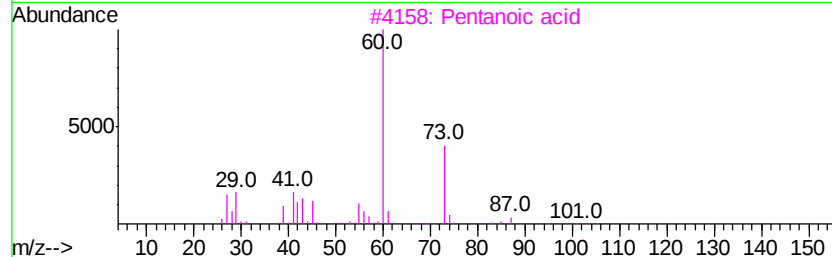
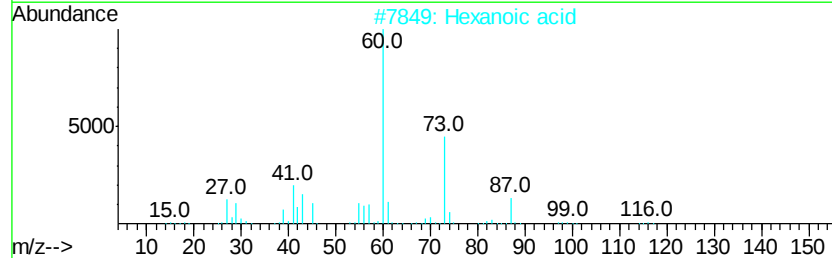
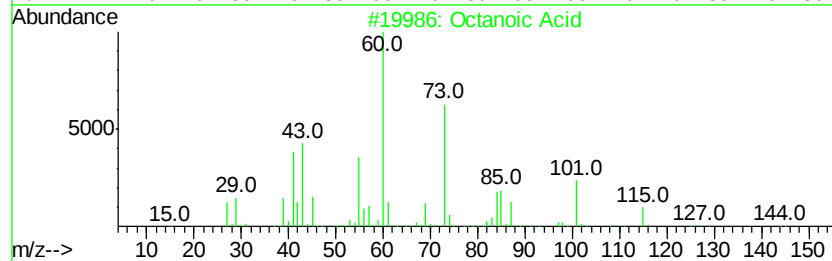
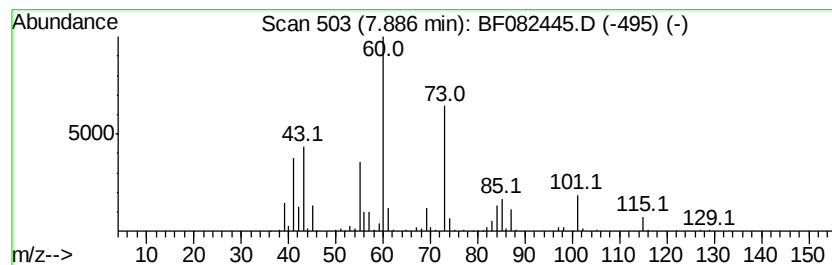
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Octanoic Acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.89	33.81 ng	1278160	Naphthalene-d8	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic Acid	144	C8H16O2	000124-07-2	91
2		Hexanoic acid	116	C6H12O2	000142-62-1	68
3		Pentanoic acid	102	C5H10O2	000109-52-4	53
4		Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	50
5		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	42



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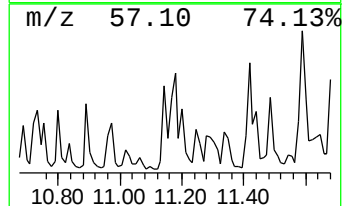
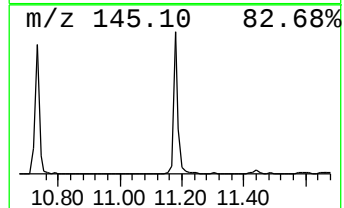
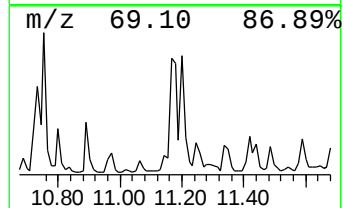
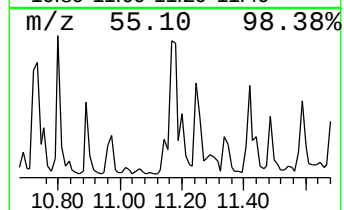
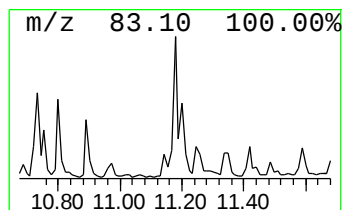
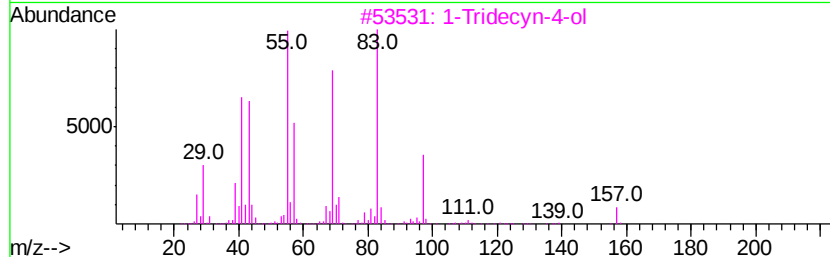
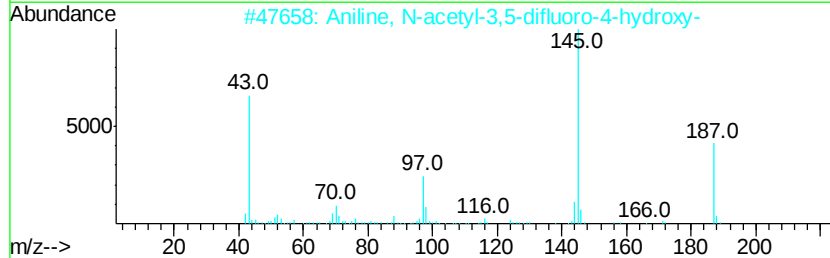
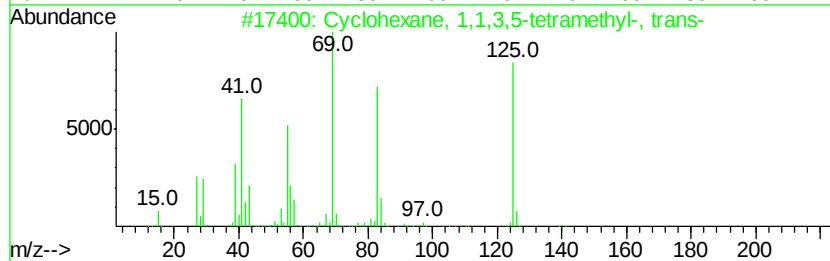
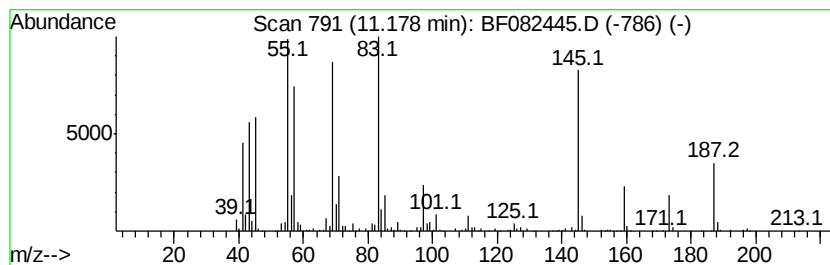
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown11.18 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	35.09 ng	2763460	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	22
2		Aniline, N-acetyl-3,5-difluoro-4...	187	C8H7F2NO2	151414-41-4	22
3		1-Tridecyn-4-ol	196	C13H24O	074646-37-0	16
4		1-Undecene, 8-methyl-	168	C12H24	074630-40-3	12
5		1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	11



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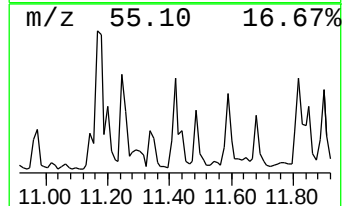
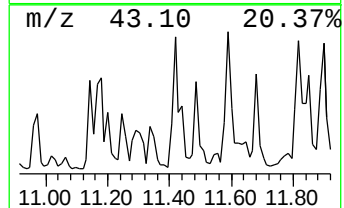
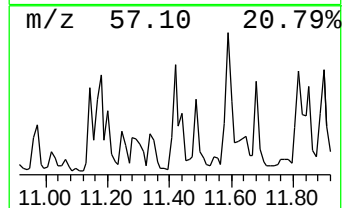
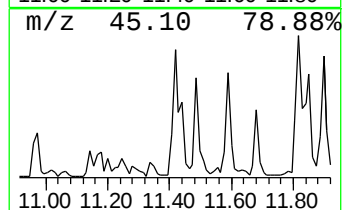
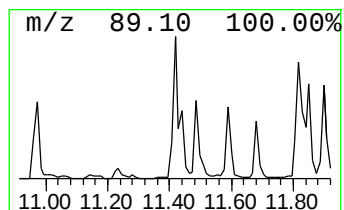
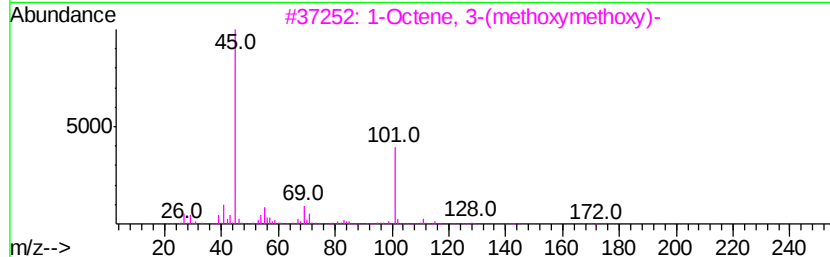
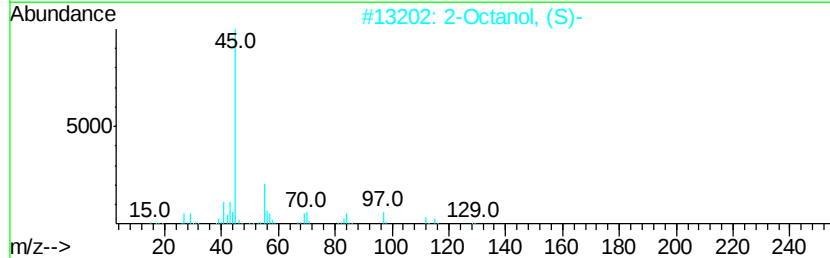
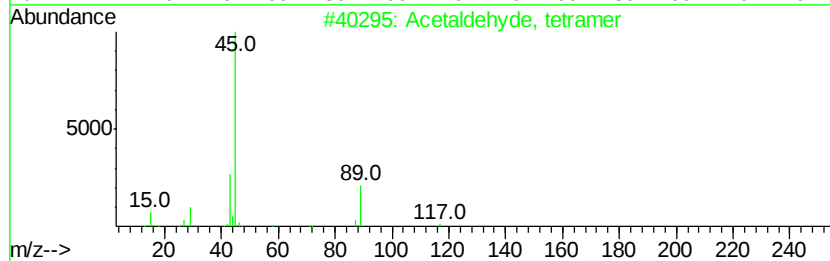
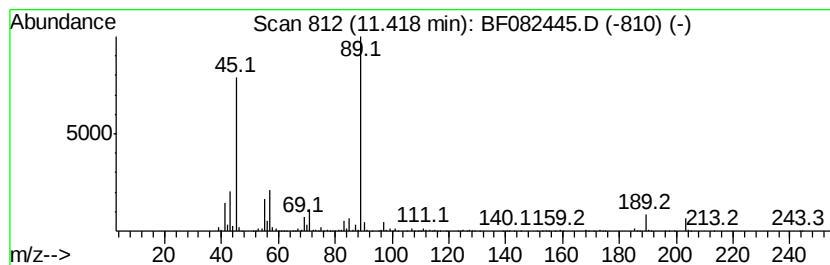
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Peak Number 3 unknown11.42 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	41.08 ng	3235340	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetaldehyde, tetramer	176	C8H16O4	000108-62-3	43
2		2-Octanol, (S)-	130	C8H18O	006169-06-8	10
3		1-Octene, 3-(methoxymethoxy)-	172	C10H20O2	088738-34-5	9
4		2-Heptanol, 4-methyl-	130	C8H18O	056298-90-9	9
5		Propanoic acid, 2-hydroxy-, buty...	146	C7H14O3	000138-22-7	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
Data File : BF082445.D
Acq On : 22 Oct 2015 11:57
Operator : UM/IZ
Sample : G4111-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

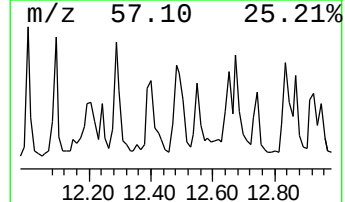
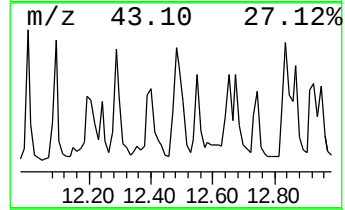
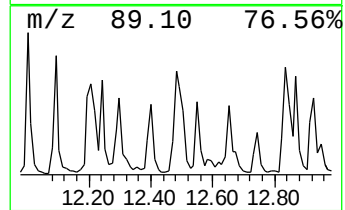
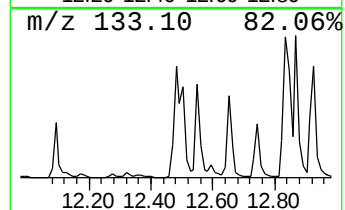
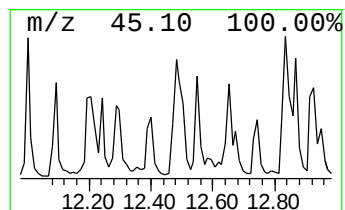
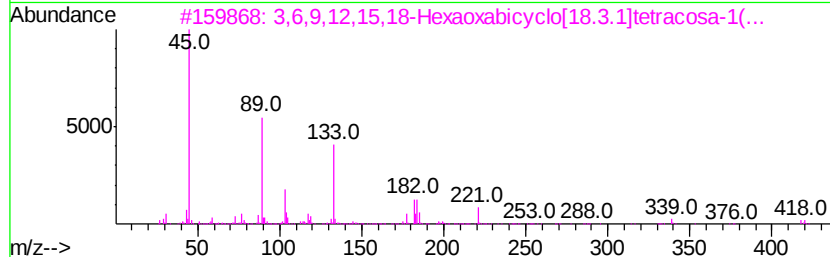
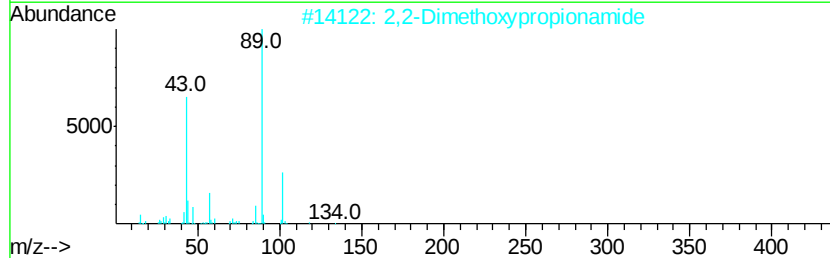
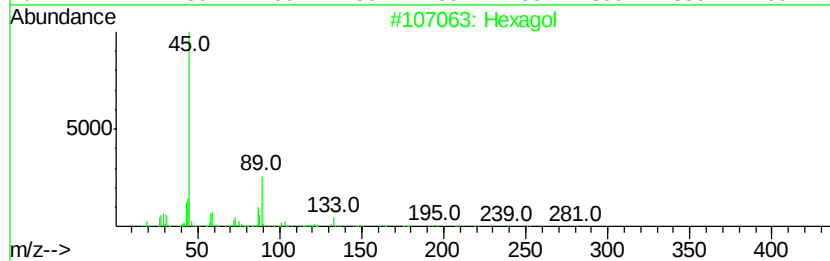
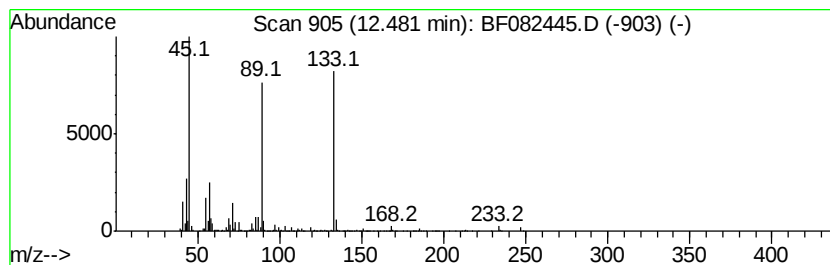
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexagol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.48	43.66 ng	3439030	Phenanthrene-d10	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexagol	282	C12H26O7	002615-15-8	74
2	2,2-Dimethoxypropionamide	133	C5H11NO3	1000237-69-7	50
3	3,6,9,12,15,18-Hexaoxabicyclo[18...]	418	C18H27BrO6	111982-22-0	40
4	Cyclobutane, 1-bromo-3-chloro-	168	C4H6BrCl	004935-03-9	38
5	Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	33



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
Data File : BF082445.D
Acq On : 22 Oct 2015 11:57
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Sample : G4111-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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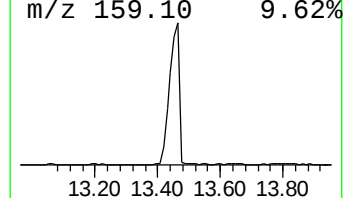
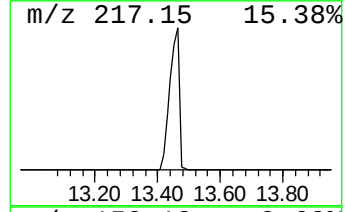
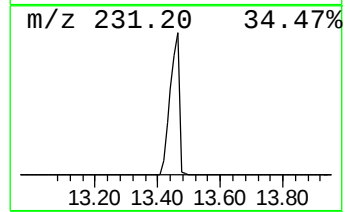
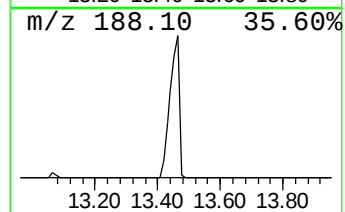
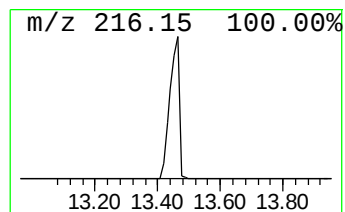
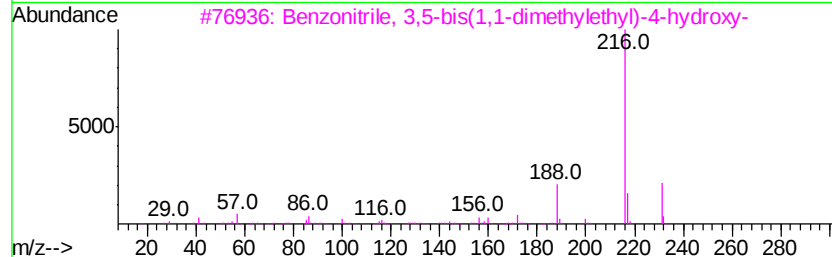
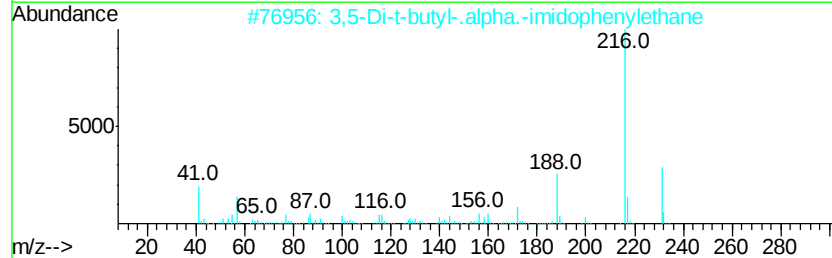
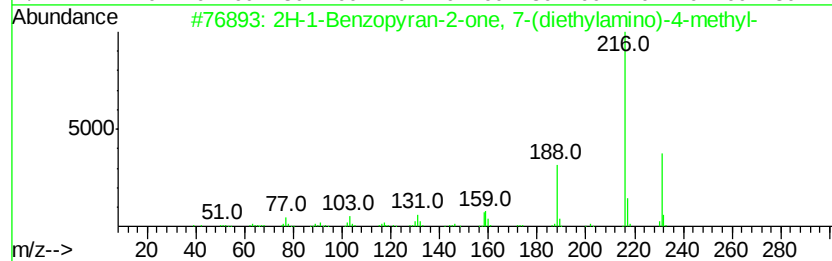
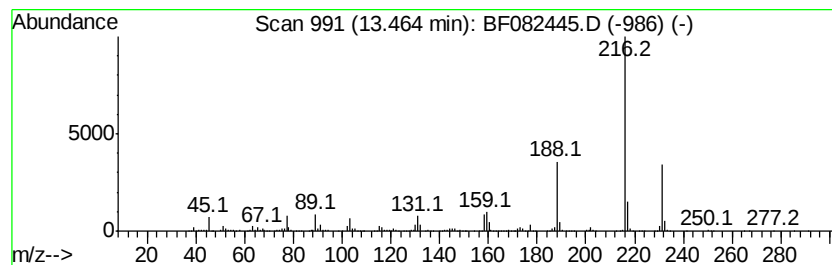
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2H-1-Benzopyran-2-one, 7-(d... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	212.30 ng	9412230	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1-Benzopyran-2-one, 7-(diethy...	231	C14H17N02	000091-44-1	98
2		3,5-Di-t-butyl-.alpha.-imidophen...	231	C16H25N	1000130-25-8	87
3		Benzonitrile, 3,5-bis(1,1-dimeth...	231	C15H21N0	001988-88-1	72
4		Phenol, 2,6-bis(1,1-dimethylethy...	231	C15H21N0	020600-84-4	72
5		2-Amino-6-(1,1-ethylenedioxyethy...	231	C12H13N3O2	089169-94-8	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
Data File : BF082445.D
Acq On : 22 Oct 2015 11:57
Operator : UM/IZ
Sample : G4111-03
Misc :
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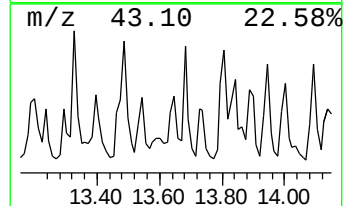
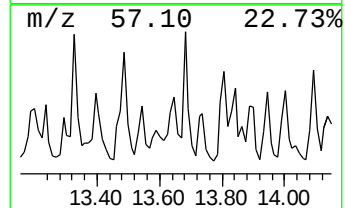
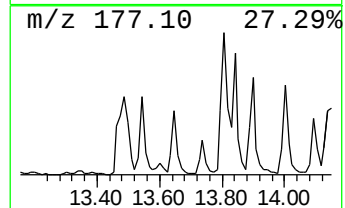
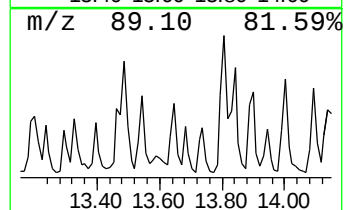
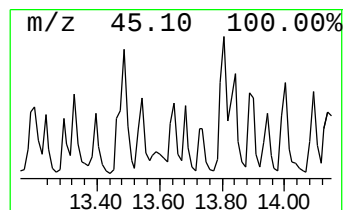
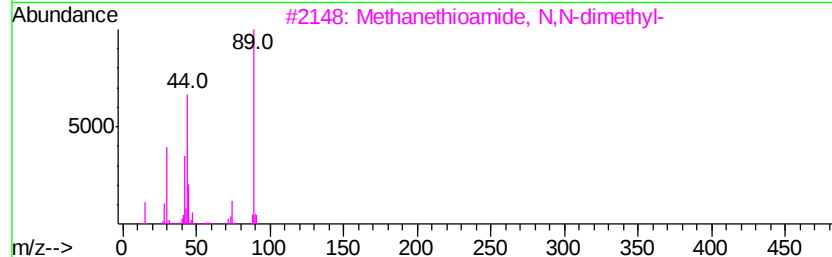
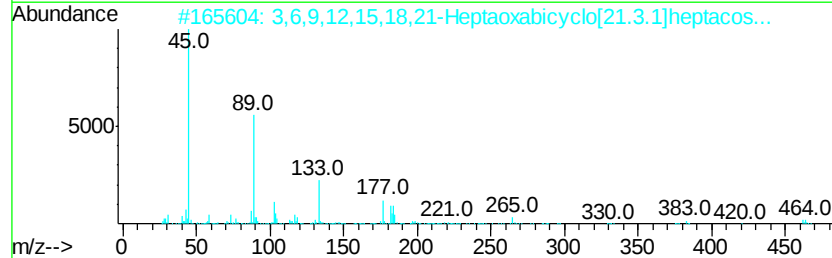
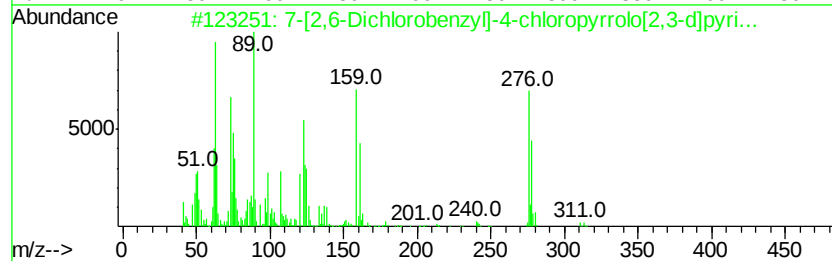
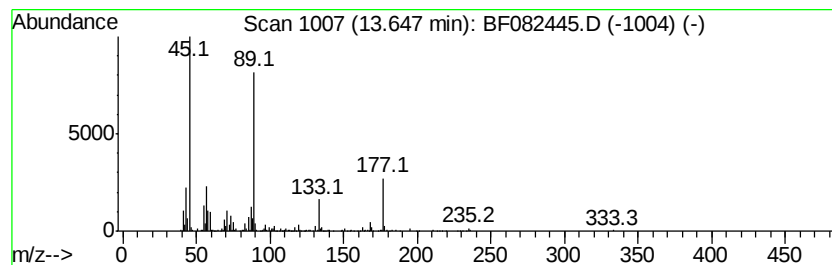
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 7-[2,6-Dichlorobenzyl]-4-ch... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.65	34.76 ng	1541080	Chrysene-d12	14.04		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-[2,6-Dichlorobenzyl]-4-chlorop...	311	C13H8Cl3N3	1000212-65-9	50	
2	3,6,9,12,15,18,21-Heptaoxabicycl...	462	C20H31BrO7	111982-23-1	43	
3	Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	42	
4	Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	42	
5	Cyclobutane, 1-bromo-3-chloro-	168	C4H6BrCl	004935-03-9	40	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
Data File : BF082445.D
Acq On : 22 Oct 2015 11:57
Operator : UM/IZ
Sample : G4111-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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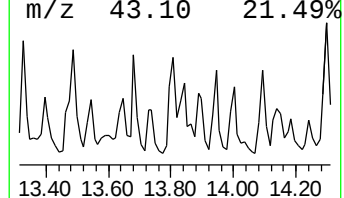
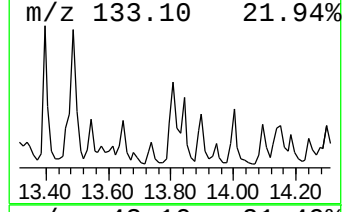
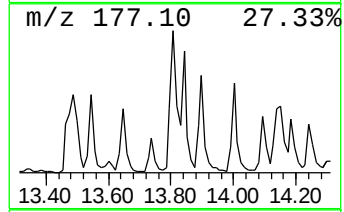
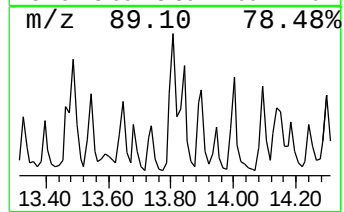
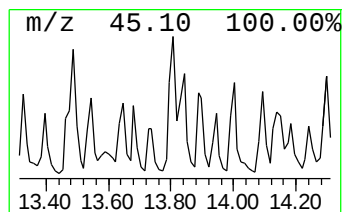
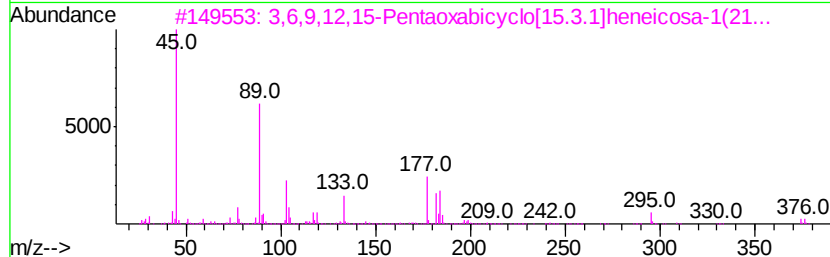
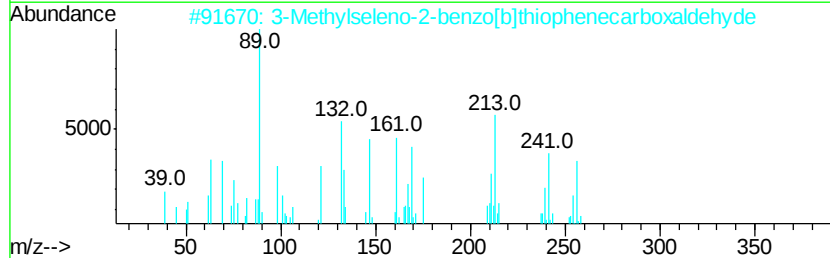
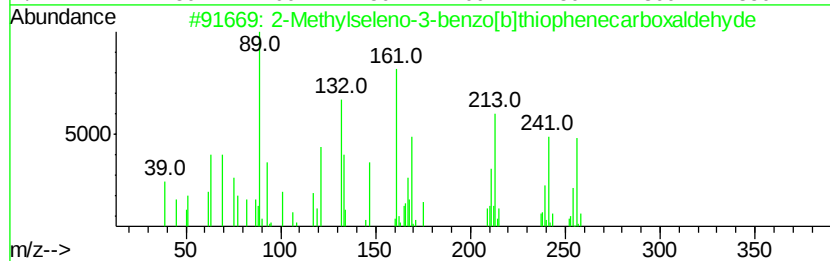
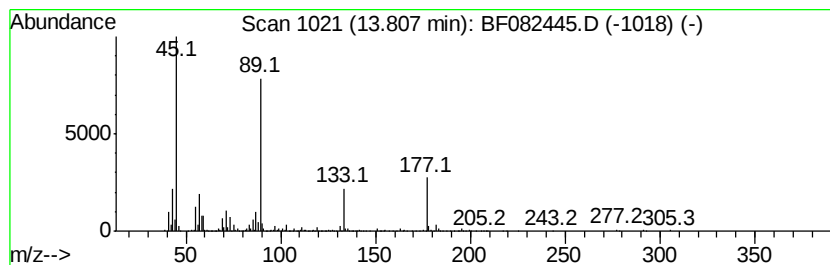
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Methylseleno-3-benzo[b]th... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	57.31 ng	2540840	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylseleno-3-benzo[b]thiophe...	256	C10H8OSse	039857-07-3	53
2		3-Methylseleno-2-benzo[b]thiophe...	256	C10H8OSse	039857-04-0	53
3		3,6,9,12,15-Pentaoxabicyclo[15.3...	374	C16H23BrO5	055440-88-5	43
4		Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	42
5		Propane, 2,2-bis(methylthio)-	136	C5H12S2	006156-18-9	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102115\
Data File : BF082445.D
Acq On : 22 Oct 2015 11:57
Operator : UM/IZ
Sample : G4111-03
Misc :
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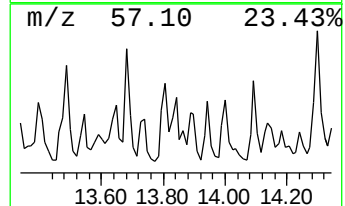
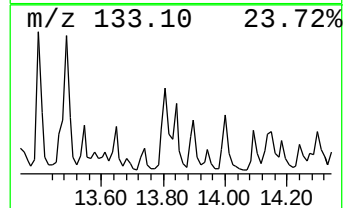
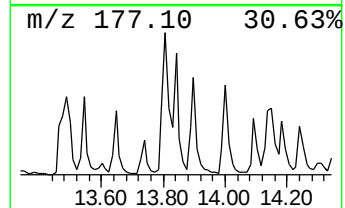
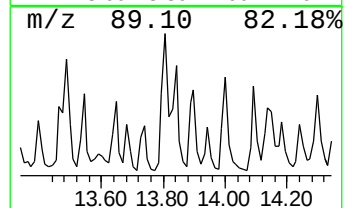
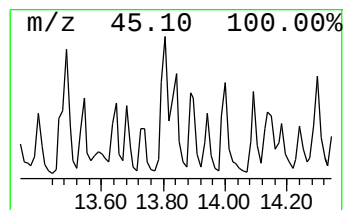
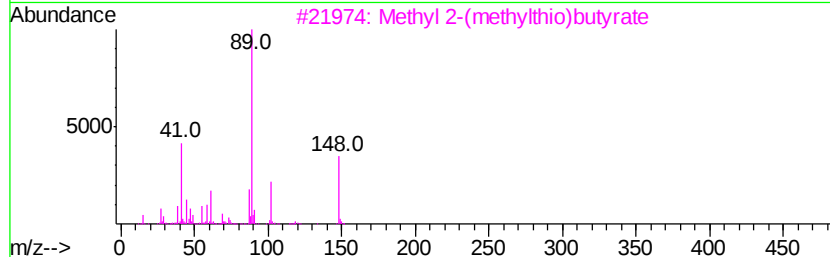
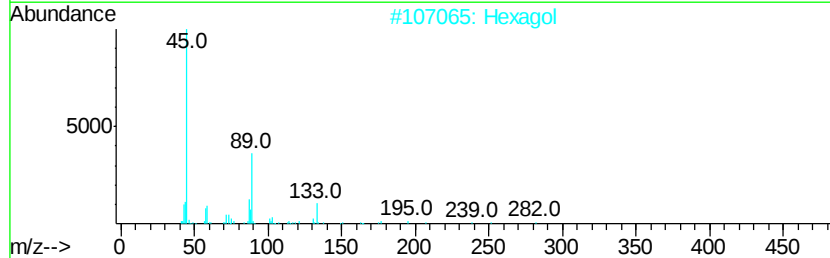
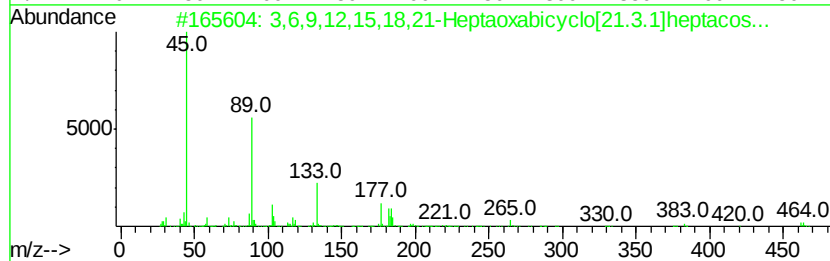
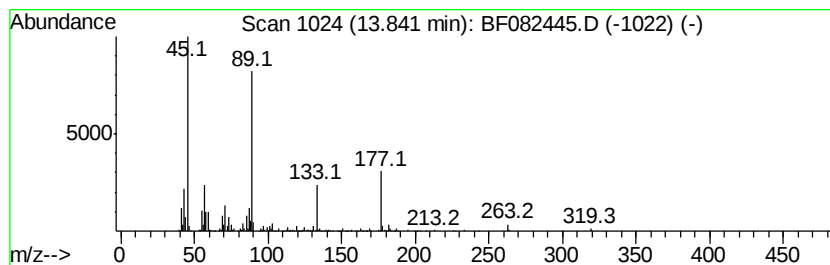
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 3,6,9,12,15,18,21-Heptaoxab... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.84	43.42 ng	1924930	Chrysene-d12	14.04		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6,9,12,15,18,21-Heptaoxabicyclo[21.3.1]heptacos...	462	C20H31BrO7	111982-23-1	64	
2	Hexagol	282	C12H26O7	002615-15-8	64	
3	Methyl 2-(methylthio)butyrate	148	C6H12O2S	051534-66-8	53	
4	2-Methylseleno-3-benzo[b]thiophe...	256	C10H8OSse	039857-07-3	53	
5	3-Methylseleno-2-benzo[b]thiophe...	256	C10H8OSse	039857-04-0	53	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
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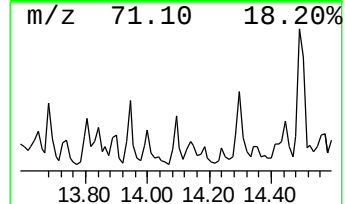
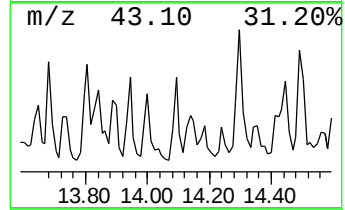
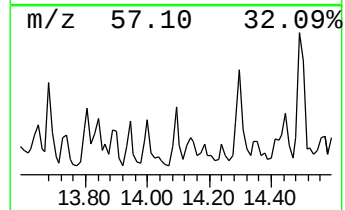
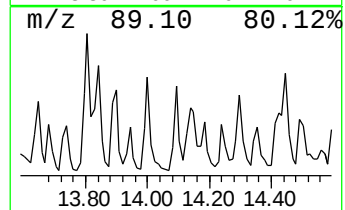
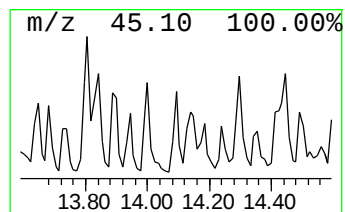
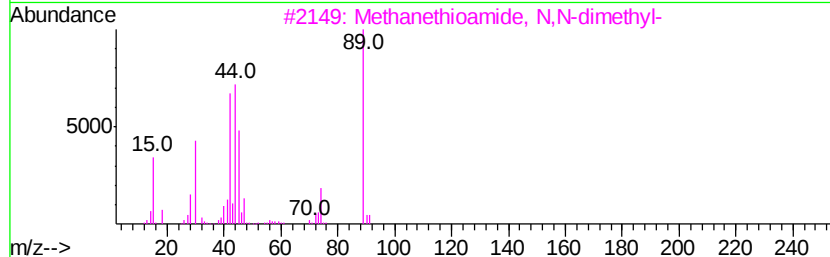
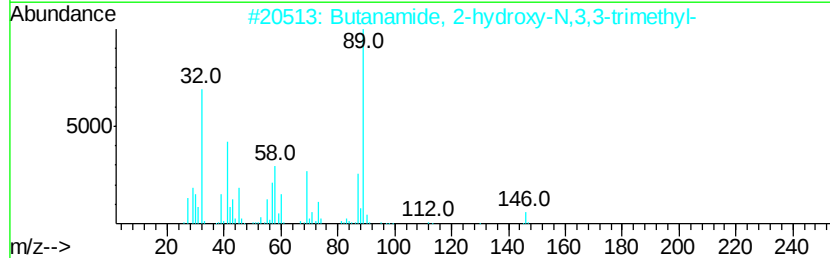
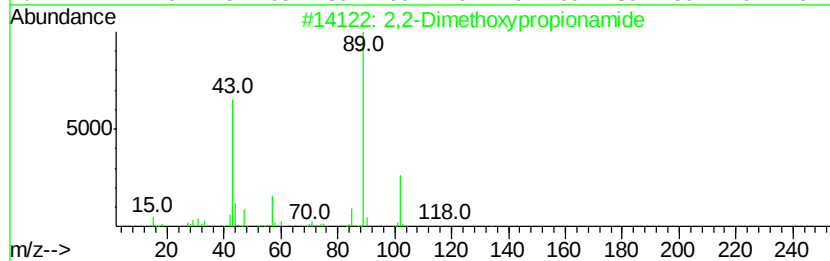
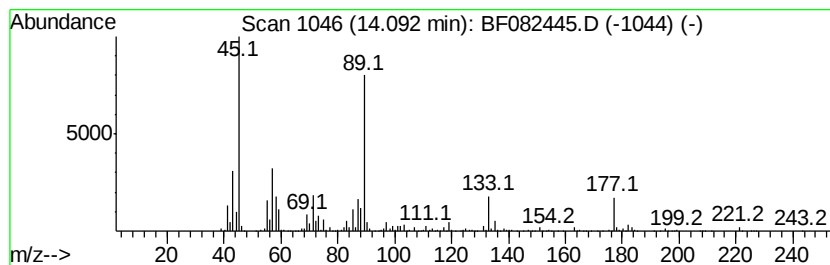
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2,2-Dimethoxypropionamide Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.09	37.08 ng	1643750	Chrysene-d12	14.04		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,2-Dimethoxypropionamide	133	C5H11NO3	1000237-69-7	53
2		Butanamide, 2-hydroxy-N,3,3-trim...	145	C7H15NO2	087919-98-0	53
3		Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	53
4		1,3-Heptadiyne, 6-(2-methoxyvetho...	268	C14H24O3Si	1000156-78-0	53
5		Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102115\
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Sample : G4111-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

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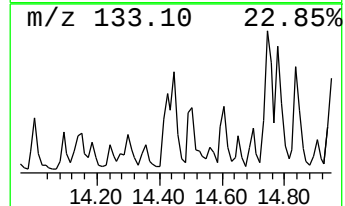
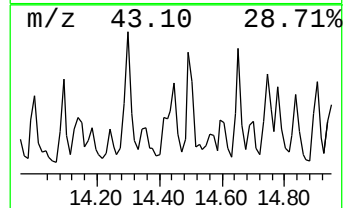
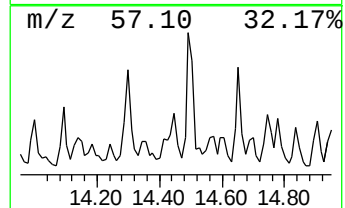
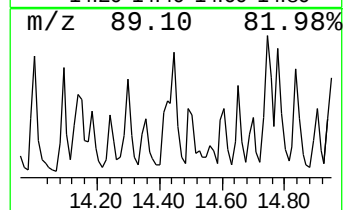
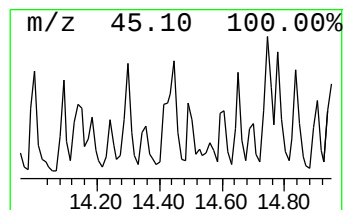
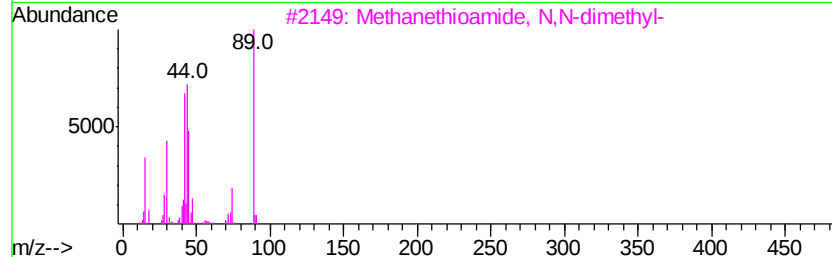
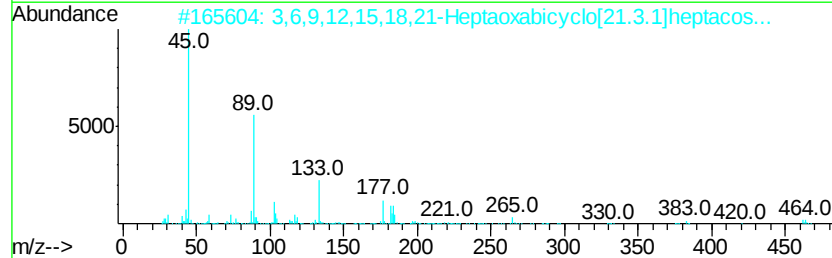
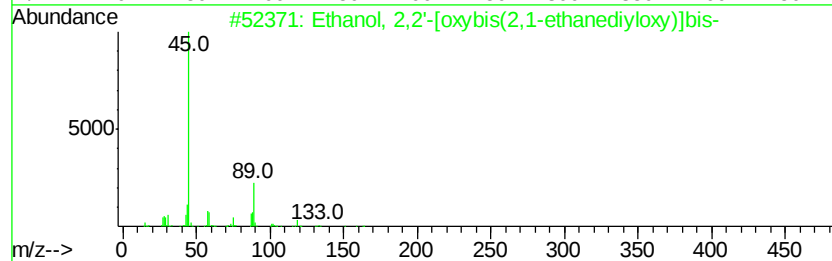
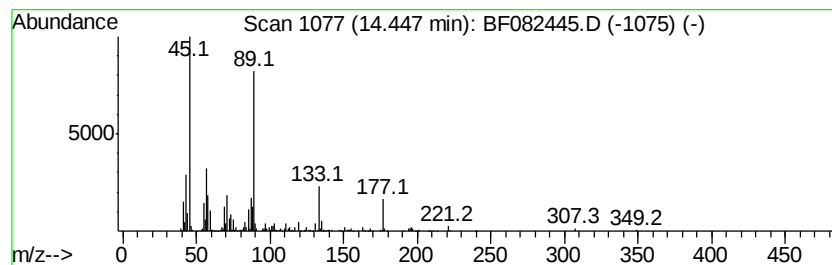
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Ethanol, 2,2'-[oxybis(2,1-e... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	40.51 ng	1796140	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2,2'-[oxybis(2,1-ethane...	194	C8H18O5	000112-60-7	72
2		3,6,9,12,15,18,21-Heptaoxabicycl...	462	C20H31BrO7	111982-23-1	64
3		Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	53
4		Butanamide, 2-hydroxy-N,3,3-trim...	145	C7H15NO2	087919-98-0	53
5		2-Methylseleno-3-benzo[b]thiophe...	256	C10H8OSse	039857-07-3	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102115\
 Data File : BF082445.D
 Acq On : 22 Oct 2015 11:57
 Operator : UM/IZ
 Sample : G4111-03
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CLEARWELL

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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
Octanoic Acid	7.89	33.8	ng	1278160	2	8.06	756144	20.0
unknown11.18	11.18	35.1	ng	2763460	4	11.30	1575240	20.0
unknown11.42	11.42	41.1	ng	3235340	4	11.30	1575240	20.0
Hexagol	12.48	43.7	ng	3439030	4	11.30	1575240	20.0
2H-1-Benzopyran-2...	13.46	212.3	ng	9412230	5	14.04	886686	20.0
7-[2,6-Dichlorobe...	13.65	34.8	ng	1541080	5	14.04	886686	20.0
2-Methylseleno-3-...	13.81	57.3	ng	2540840	5	14.04	886686	20.0
3,6,9,12,15,18,21...	13.84	43.4	ng	1924930	5	14.04	886686	20.0
2,2-Dimethoxyprop...	14.09	37.1	ng	1643750	5	14.04	886686	20.0
Ethanol, 2,2'-[ox...	14.45	40.5	ng	1796140	5	14.04	886686	20.0