

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
 Data File : BF133047.D
 Acq On : 26 Apr 2023 01:34
 Operator : CG\JU
 Sample : 02390-11 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SPC-3WC-23-04-17

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133047.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.281	29	33	37	rVB2	59401	80589	2.72%	0.169%
2	3.710	269	276	284	rVB	116191	167623	5.66%	0.351%
3	4.416	392	396	400	rBV	63797	74322	2.51%	0.156%
4	4.916	477	481	488	rVB	82170	96834	3.27%	0.203%
5	5.010	493	497	501	rVB	63680	60682	2.05%	0.127%
6	5.199	526	529	532	rBV	125078	107772	3.64%	0.226%
7	5.310	544	548	550	rBV	181328	176968	5.98%	0.371%
8	5.340	550	553	557	rVB	1664351	1524865	51.53%	3.194%
9	5.575	586	593	601	rVB	70368	84500	2.86%	0.177%
10	6.263	708	710	714	rBV	58465	51339	1.73%	0.108%
11	6.340	718	723	726	rBV4	52431	58860	1.99%	0.123%
12	6.375	726	729	731	rBV	1792211	1552564	52.47%	3.252%
13	6.428	731	738	739	rBV	359447	801613	27.09%	1.679%
14	6.446	739	741	742	rBV	294604	217474	7.35%	0.456%
15	6.569	759	762	763	rBV	135795	116601	3.94%	0.244%
16	6.740	788	791	794	rBV	1757287	1362104	46.03%	2.853%
17	6.910	815	820	824	rVB	154658	134342	4.54%	0.281%
18	7.063	843	846	849	rBV	55850	54607	1.85%	0.114%
19	7.140	856	859	864	rBV4	47486	73628	2.49%	0.154%
20	7.187	864	867	873	rBV2	92798	208400	7.04%	0.437%
21	7.298	880	886	888	rBV	1420868	1090689	36.86%	2.285%
22	7.316	888	889	892	rVV	122876	72225	2.44%	0.151%
23	7.345	892	894	897	rVB	71133	55400	1.87%	0.116%
24	7.422	904	907	910	rBV	609222	455177	15.38%	0.953%
25	7.822	972	975	978	rVB2	107622	78254	2.64%	0.164%
26	7.869	979	983	986	rBV	169764	169956	5.74%	0.356%
27	7.963	997	999	1002	rVB	148509	118053	3.99%	0.247%
28	8.022	1006	1009	1011	rBV	1983626	1678134	56.71%	3.515%
29	8.045	1011	1013	1016	rVB	475022	332858	11.25%	0.697%
30	8.204	1032	1040	1048	rVB3	123907	152213	5.14%	0.319%
31	8.734	1128	1130	1134	rVB	166227	122981	4.16%	0.258%
32	8.781	1134	1138	1144	rBV5	30859	58026	1.96%	0.122%
33	8.834	1144	1147	1151	rVV	102351	76718	2.59%	0.161%
34	9.051	1181	1184	1186	rBV2	65040	60935	2.06%	0.128%
35	9.098	1189	1192	1195	rVB	2408257	1788272	60.43%	3.746%
36	9.198	1204	1209	1212	rBV2	172748	183414	6.20%	0.384%

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Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	9.298	1223	1226	1229	rBV	185357	166508	5.63%	0.349%
38	9.357	1233	1236	1240	rVB	62309	76565	2.59%	0.160%
39	9.434	1246	1249	1252	rBV2	114496	115415	3.90%	0.242%
40	9.551	1266	1269	1272	rBV	341913	234124	7.91%	0.490%
41	9.622	1278	1281	1288	rBV2	82697	125395	4.24%	0.263%
42	9.775	1301	1307	1310	rVV	2354498	1865295	63.04%	3.907%
43	9.898	1325	1328	1329	rBV	121274	75076	2.54%	0.157%
44	9.916	1329	1331	1335	rVV	128375	118036	3.99%	0.247%
45	9.957	1335	1338	1340	rVV	640586	529677	17.90%	1.109%
46	9.975	1340	1341	1344	rVB	107679	72152	2.44%	0.151%
47	10.016	1344	1348	1351	rVB3	159928	177193	5.99%	0.371%
48	10.057	1353	1355	1360	rBV	73108	72690	2.46%	0.152%
49	10.116	1363	1365	1370	rBV3	43341	55196	1.87%	0.116%
50	10.186	1374	1377	1380	rBV4	36238	52152	1.76%	0.109%
51	10.286	1390	1394	1397	rBV	743390	589067	19.91%	1.234%
52	10.322	1397	1400	1406	rVB3	232364	250375	8.46%	0.524%
53	10.428	1416	1418	1422	rVV	845606	697975	23.59%	1.462%
54	10.481	1424	1427	1432	rVB2	220597	263135	8.89%	0.551%
55	10.557	1437	1440	1444	rBV2	1240457	1081826	36.56%	2.266%
56	10.604	1445	1448	1450	rVB	245103	204125	6.90%	0.428%
57	10.628	1450	1452	1459	rVB	94326	92313	3.12%	0.193%
58	10.686	1460	1462	1463	rBV2	84587	68895	2.33%	0.144%
59	10.757	1471	1474	1476	rBV	58003	69918	2.36%	0.146%
60	10.786	1476	1479	1482	rVV2	73063	82329	2.78%	0.172%
61	10.822	1482	1485	1487	rBV2	58833	69203	2.34%	0.145%
62	10.869	1490	1493	1496	rBV	351764	291243	9.84%	0.610%
63	11.057	1523	1525	1529	rBV	138877	134397	4.54%	0.282%
64	11.139	1536	1539	1542	rBV3	56923	83865	2.83%	0.176%
65	11.186	1545	1547	1550	rVB	400679	333120	11.26%	0.698%
66	11.257	1555	1559	1561	rBV	1947739	1659507	56.08%	3.476%
67	11.280	1561	1563	1565	rVB	509285	357657	12.09%	0.749%
68	11.328	1569	1571	1573	rVB	109757	82190	2.78%	0.172%
69	11.451	1590	1592	1595	rBV2	105623	84866	2.87%	0.178%
70	11.757	1641	1644	1646	rBV	121450	118464	4.00%	0.248%
71	11.792	1646	1650	1654	rVV2	498434	725609	24.52%	1.520%
72	11.875	1660	1664	1668	rVB3	255235	324196	10.96%	0.679%
73	11.928	1668	1673	1677	rBV	854384	883916	29.87%	1.851%
74	12.045	1691	1693	1696	rBV2	132137	170192	5.75%	0.356%
75	12.075	1696	1698	1701	rVV	183241	162247	5.48%	0.340%
76	12.469	1762	1765	1767	rBV	772292	657660	22.22%	1.378%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	12.692	1801	1803	1808	rVB	625324	572436	19.34%	1.199%
78	12.845	1825	1829	1835	rVB	1354670	1457523	49.26%	3.053%
79	12.910	1837	1840	1843	rVV	369108	375122	12.68%	0.786%
80	13.027	1857	1860	1864	rVB	1823203	1429078	48.29%	2.993%
81	13.122	1873	1876	1881	rVB2	776261	671489	22.69%	1.407%
82	13.657	1965	1967	1971	rVV2	247005	227277	7.68%	0.476%
83	13.692	1971	1973	1977	rVB3	223588	197593	6.68%	0.414%
84	13.892	2003	2007	2010	rBV3	1918201	2044429	69.09%	4.282%
85	13.916	2010	2011	2016	rVB	332687	257709	8.71%	0.540%
86	14.033	2028	2031	2036	rVB	994673	936814	31.66%	1.962%
87	14.110	2040	2044	2047	rBV	2995884	2959108	100.00%	6.198%
88	14.263	2067	2070	2076	rVB2	934095	844573	28.54%	1.769%
89	14.345	2082	2084	2087	rBV	232822	195384	6.60%	0.409%
90	14.427	2094	2098	2101	rVB2	1410636	1421472	48.04%	2.977%
91	14.710	2143	2146	2149	rVB	683520	653045	22.07%	1.368%
92	14.916	2178	2181	2183	rBV	354805	490344	16.57%	1.027%
93	15.151	2219	2221	2225	rVB	442895	457524	15.46%	0.958%
94	15.274	2235	2242	2243	rBV3	757086	1449051	48.97%	3.035%
95	15.321	2247	2250	2253	rVB	889489	941622	31.82%	1.972%
96	15.898	2345	2348	2352	rBV3	235944	317518	10.73%	0.665%
97	16.410	2432	2435	2442	rVB2	277685	450096	15.21%	0.943%
98	16.563	2457	2461	2467	rVB	379342	594550	20.09%	1.245%
99	16.721	2485	2488	2493	rVB	403140	645006	21.80%	1.351%
100	17.698	2648	2654	2663	rVB	605529	1379813	46.63%	2.890%

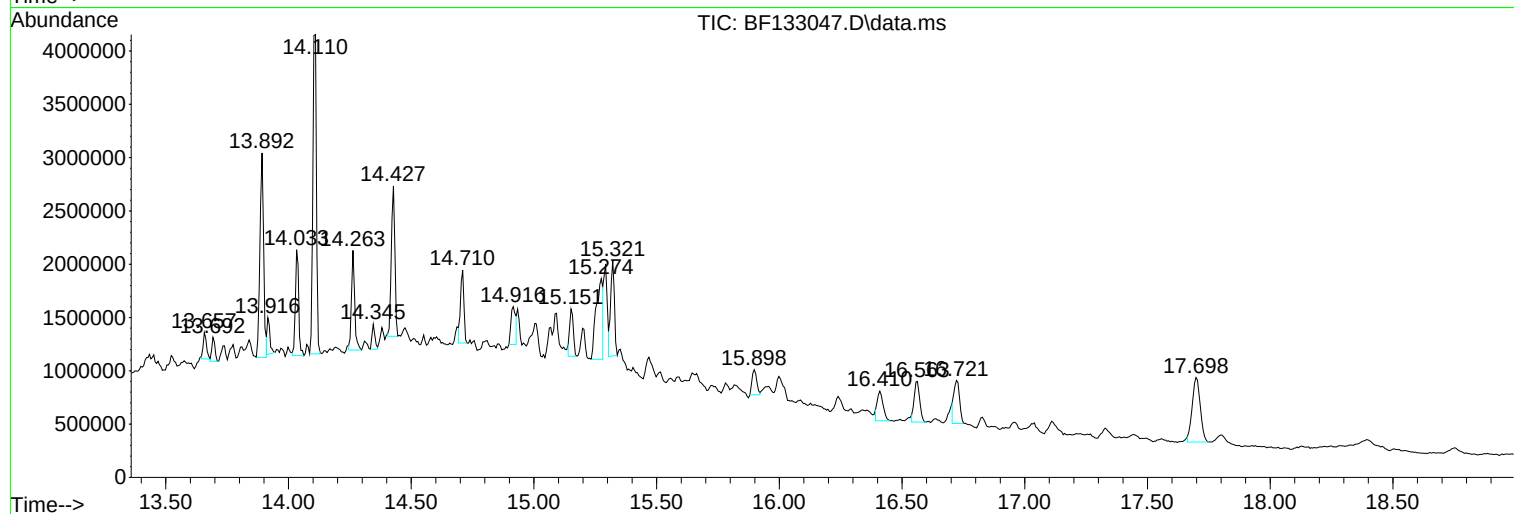
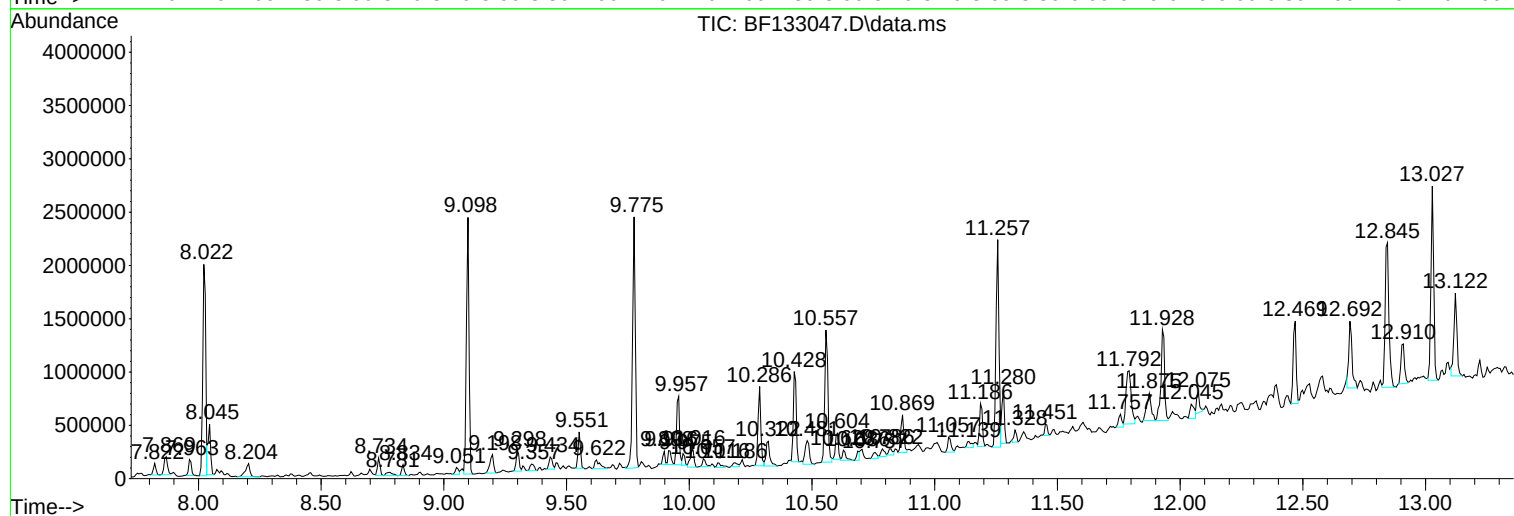
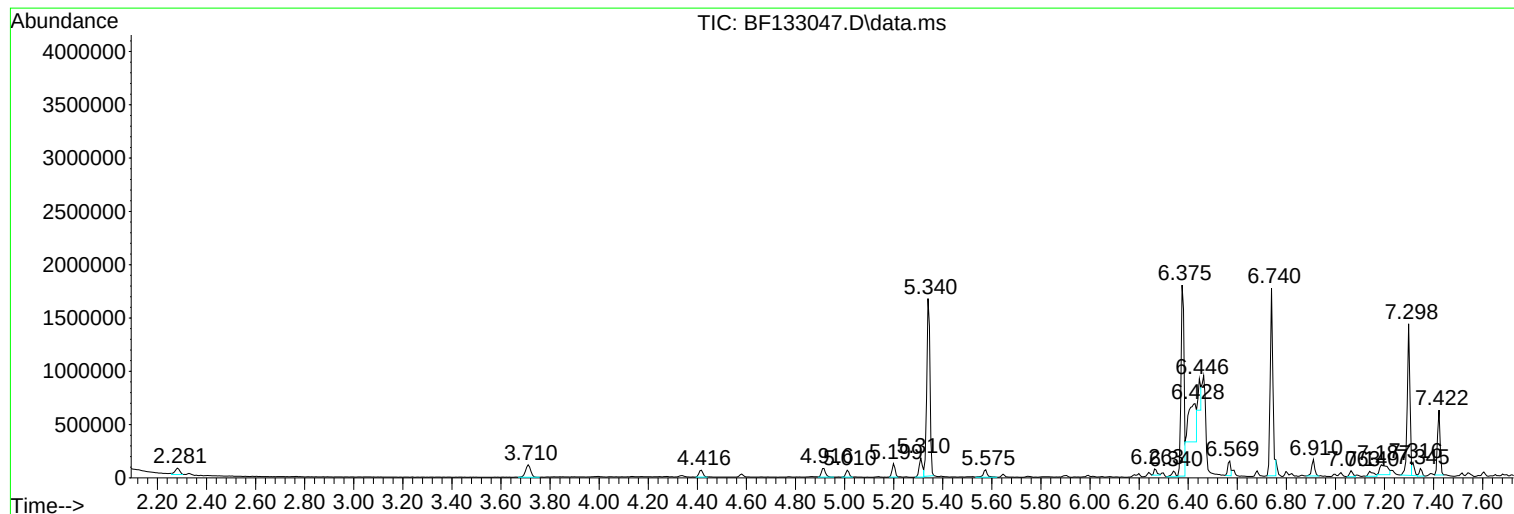
Sum of corrected areas: 47741332

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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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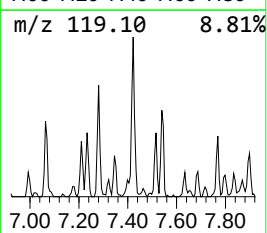
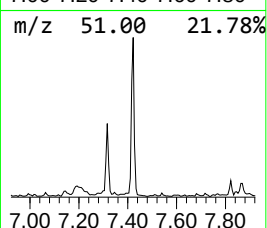
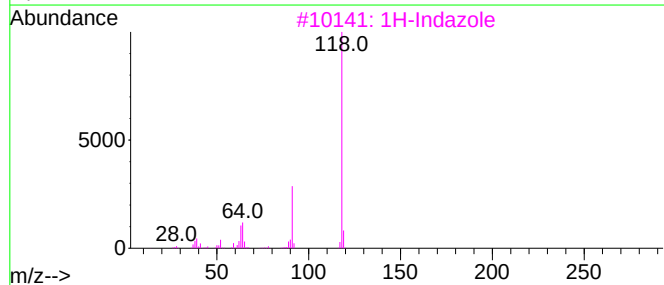
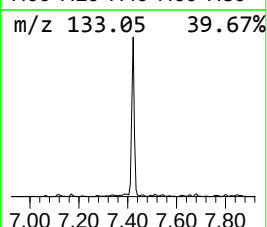
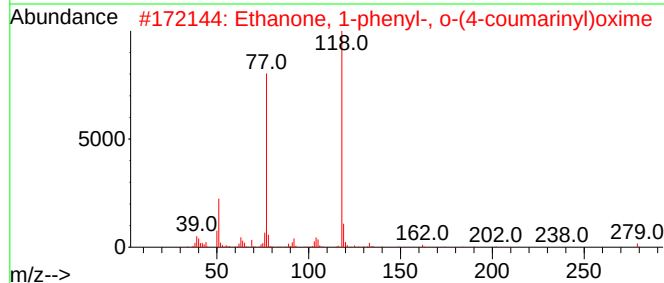
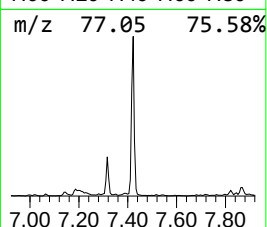
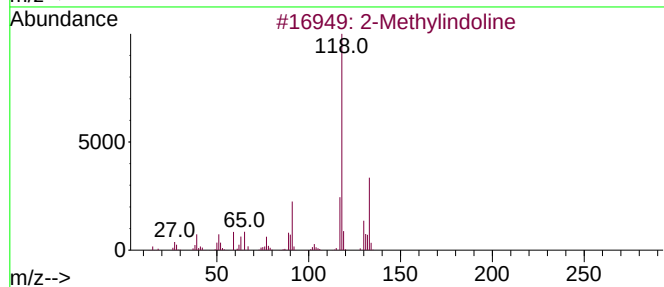
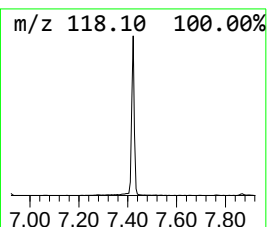
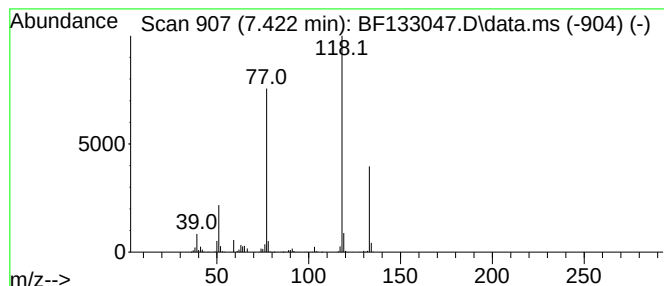
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.422 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.422	5.42 ng	455177	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylindoline	133	C9H11N	006872-06-6	43
2			Ethanone, 1-phenyl-, o-(4-coumar...	279	C17H13NO3	034605-11-3	27
3			1H-Indazole	118	C7H6N2	000271-44-3	22
4			1H-Benzimidazole	118	C7H6N2	000051-17-2	22
5			Benzofuran	118	C8H6O	000271-89-6	22



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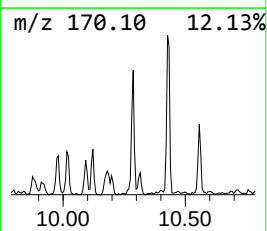
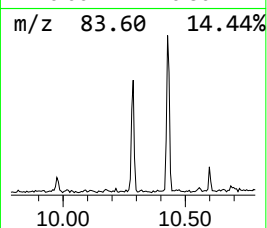
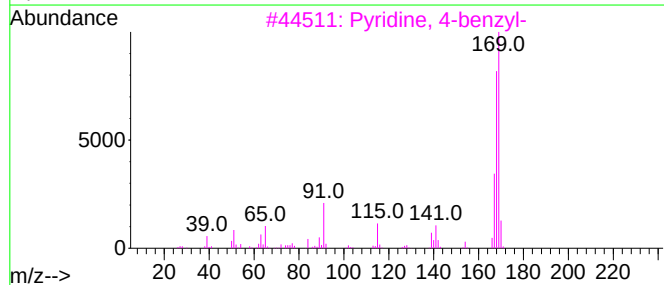
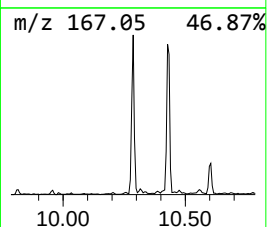
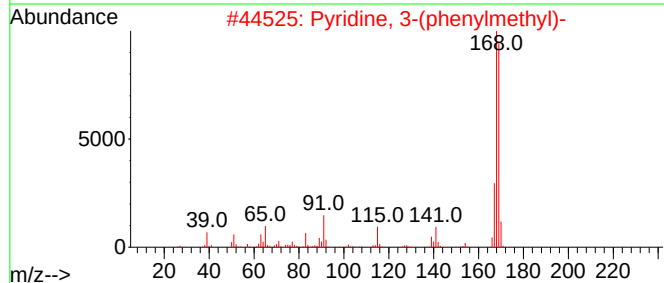
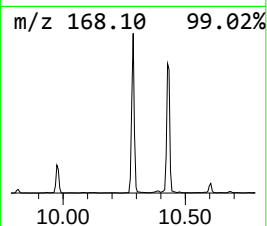
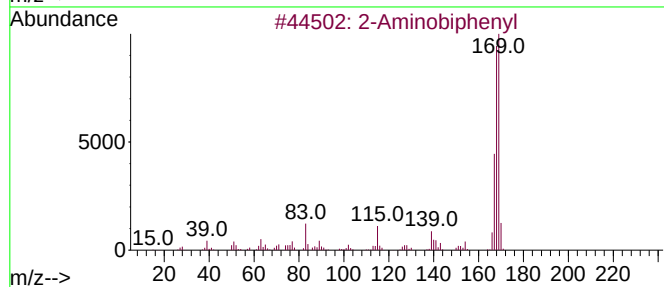
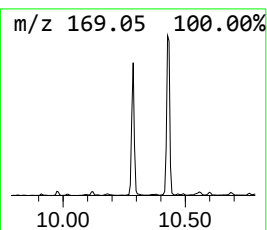
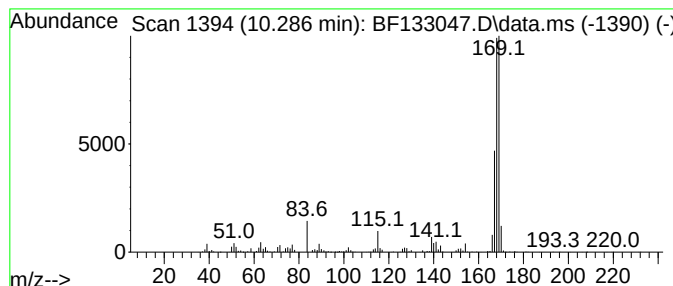
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Aminobiphenyl Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.286	6.32 ng	589067	Acenaphthene-d10	9.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Aminobiphenyl	169	C12H11N	000090-41-5	94
2			Pyridine, 3-(phenylmethyl)-	169	C12H11N	000620-95-1	74
3			Pyridine, 4-benzyl-	169	C12H11N	002116-65-6	68
4			Diphenylamine	169	C12H11N	000122-39-4	64
5			2-Methyl-3-phenylpyridine	169	C12H11N	003256-89-1	64



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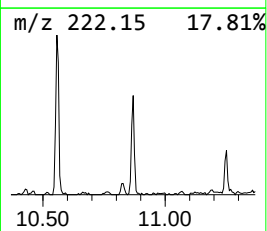
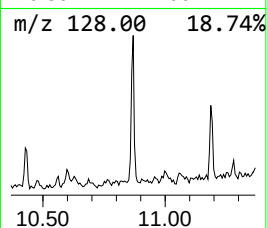
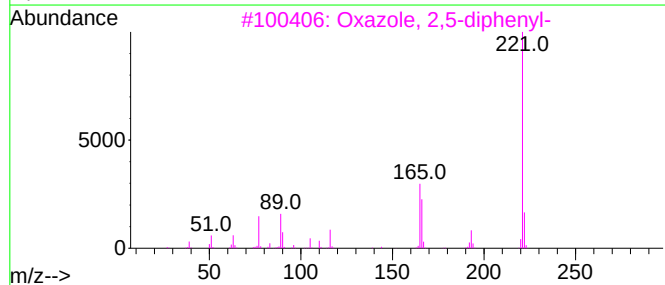
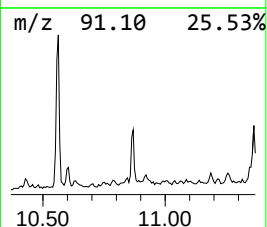
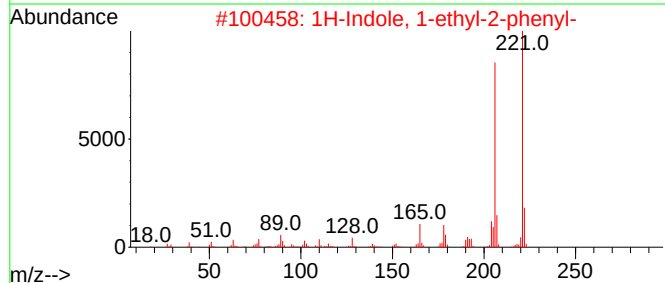
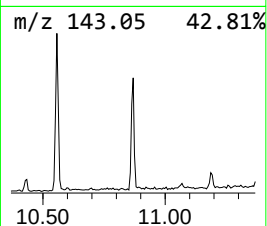
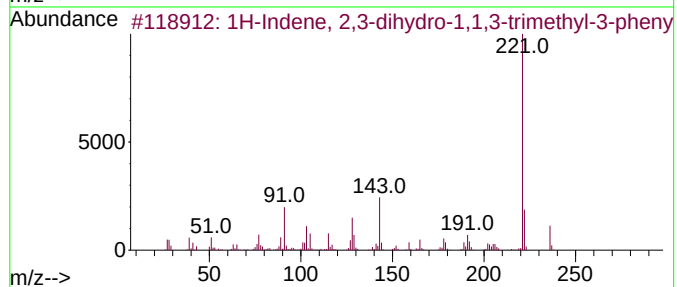
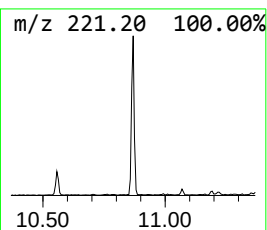
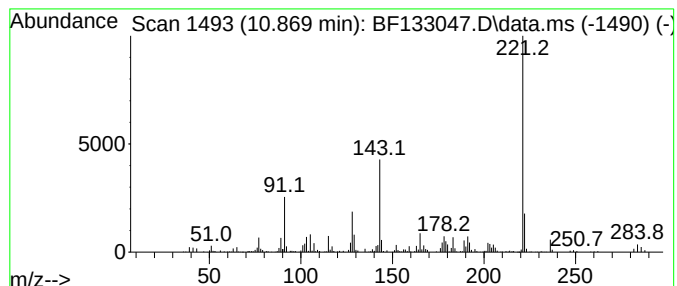
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.869	3.51 ng	291243	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	90
2			1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	49
3			Oxazole, 2,5-diphenyl-	221	C15H11NO	000092-71-7	46
4			Benzo[h]quinoline, 2,3,4-trimethyl-	221	C16H15N	063042-66-0	46
5			1H-1,2,3-Triazole, 4,5-diphenyl-	221	C14H11N3	005533-73-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
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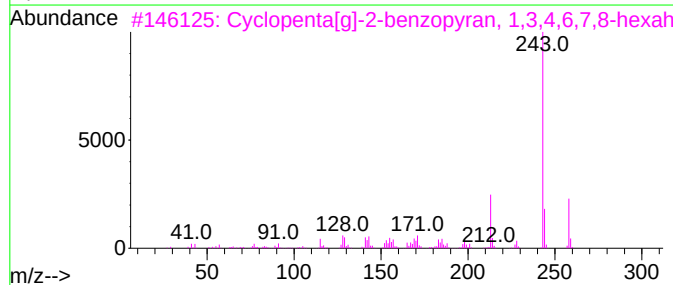
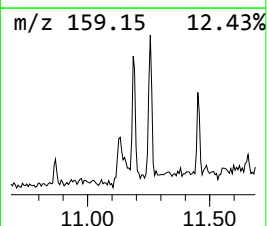
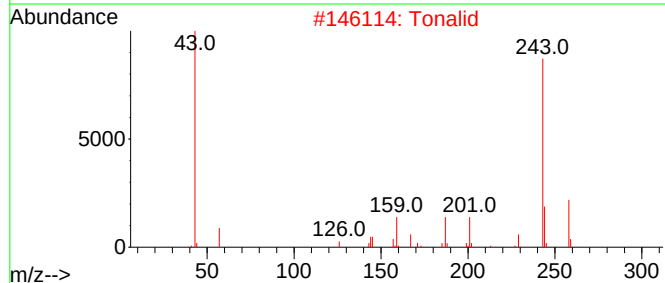
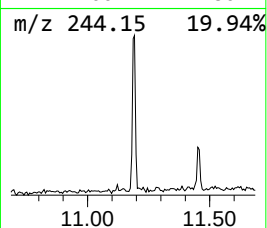
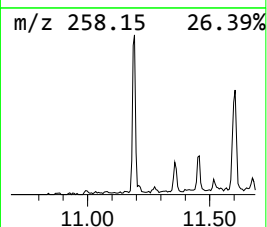
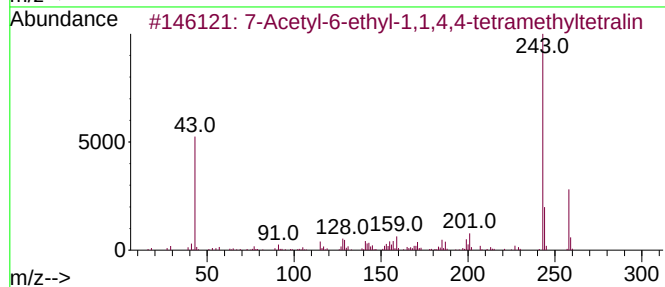
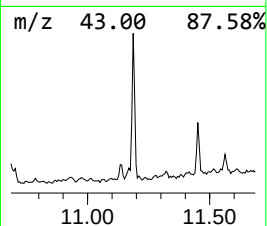
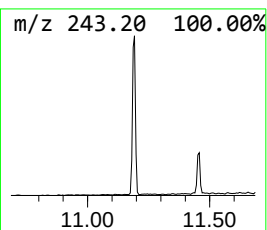
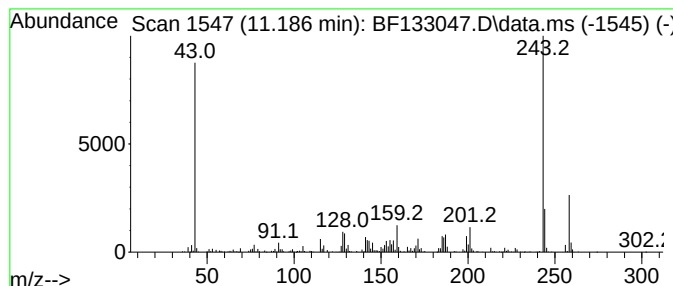
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 7-Acetyl-6-ethyl-1,1,4,4-te... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.186	4.01 ng	333120	Phenanthrene-d10	11.257	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Acetyl-6-ethyl-1,1,4,4-tetrame...	258	C18H26O	000088-29-9	98
2	Tonalid	258	C18H26O	021145-77-7	87
3	Cyclopenta[g]-2-benzopyran, 1,3,...	258	C18H26O	001222-05-5	81
4	Luvangetin	258	C15H14O4	000483-92-1	59
5	Brayelin	258	C15H14O4	006054-10-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
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ClientSampleId :
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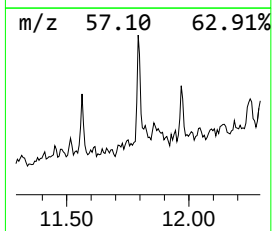
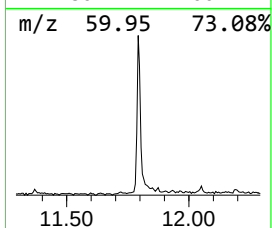
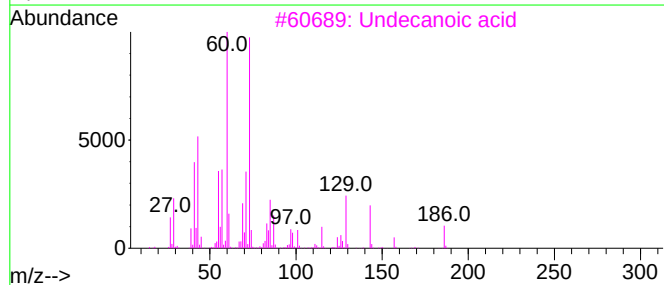
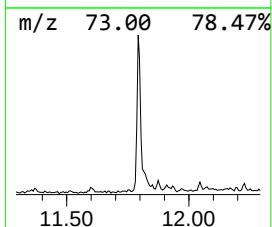
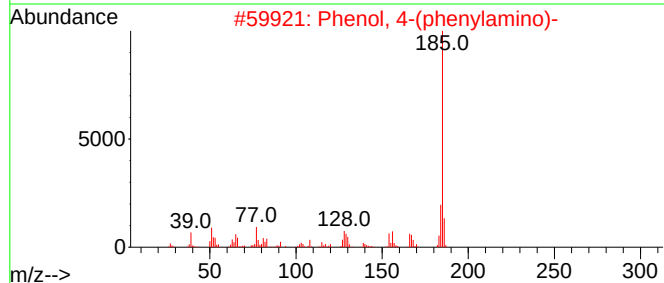
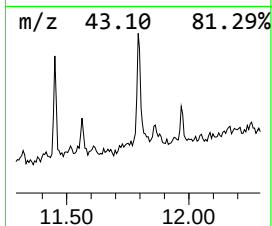
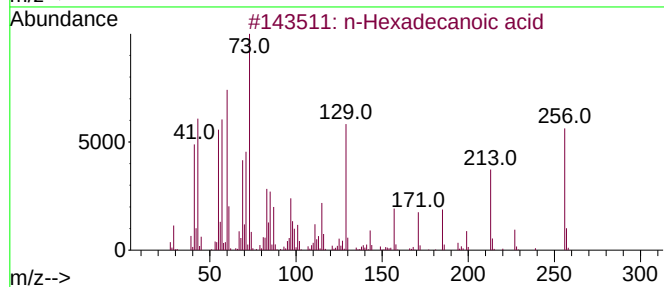
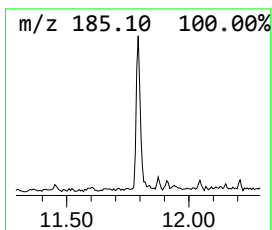
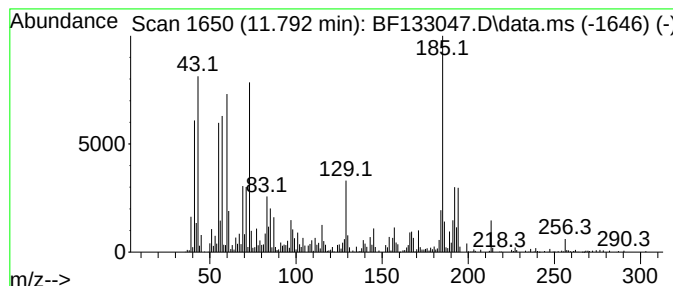
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.792	8.74 ng	725609	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2			Phenol, 4-(phenylamino)-	185	C12H11NO	000122-37-2	93
3			Undecanoic acid	186	C11H22O2	000112-37-8	62
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	50
5			3-Hydroxydiphenylamine	185	C12H11NO	000101-18-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
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Operator : CG\JU
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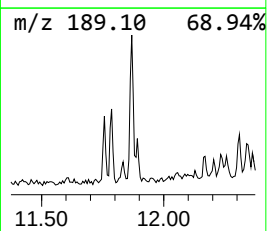
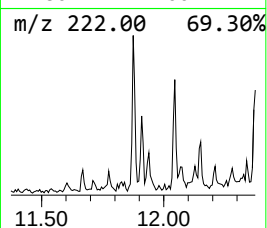
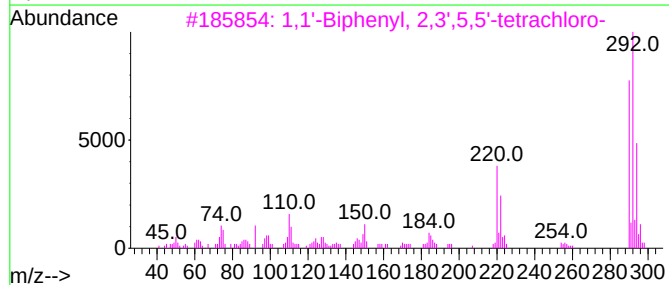
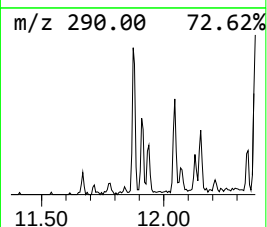
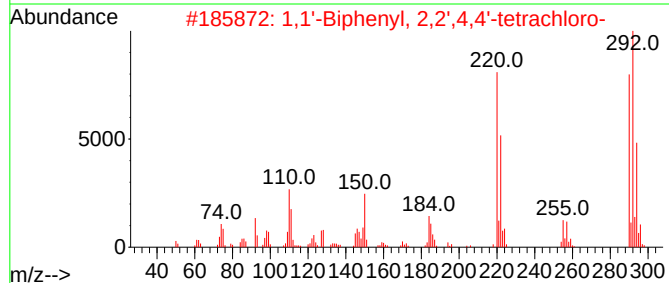
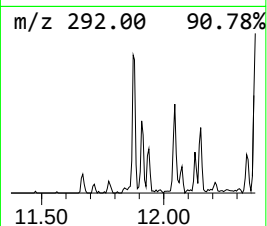
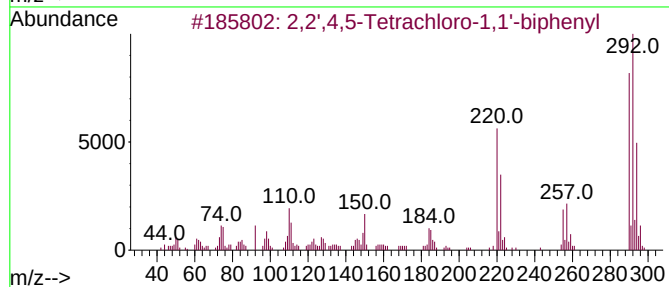
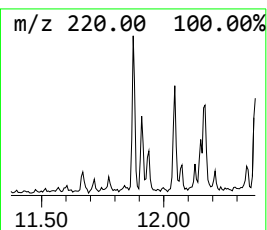
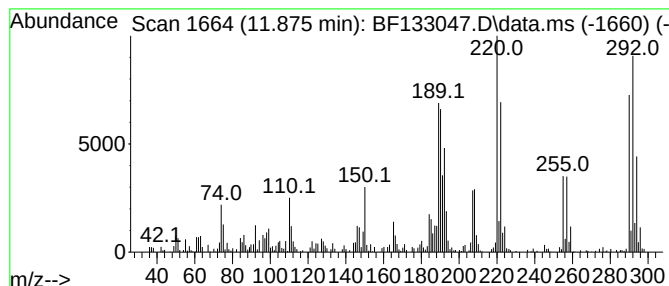
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.875	3.91 ng	324196	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99		
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
3	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99		
4	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99		
5	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
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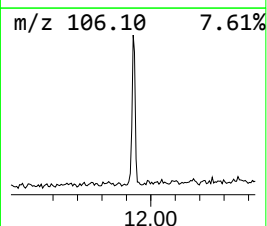
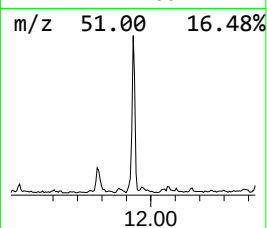
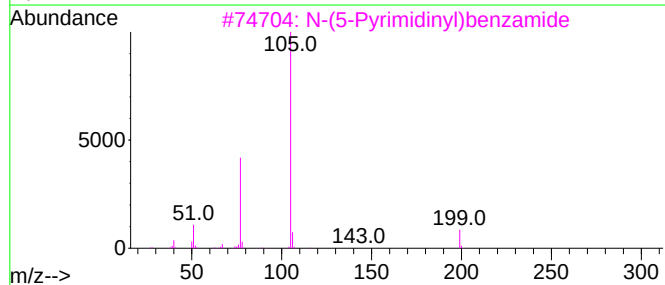
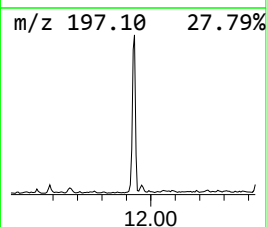
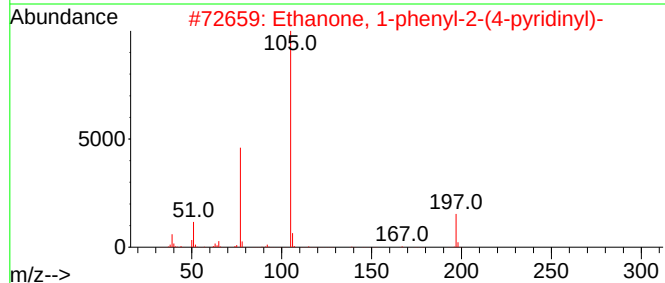
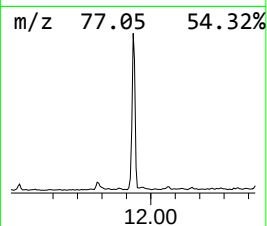
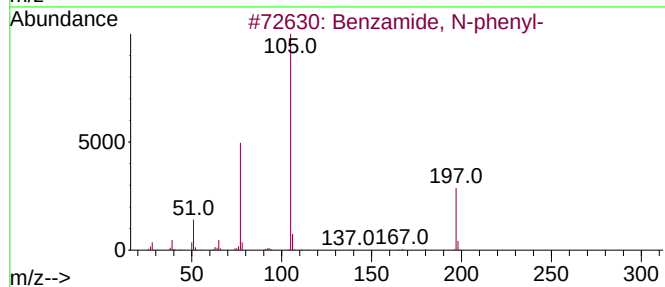
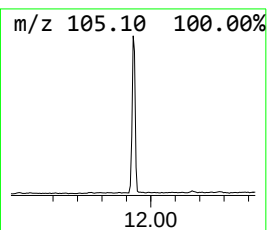
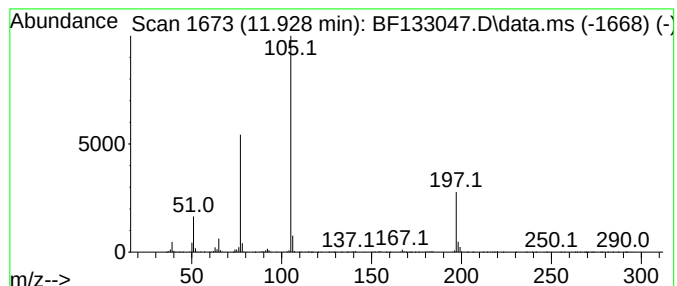
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzamide, N-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.928	10.65 ng	883916	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-phenyl-	197	C13H11NO	000093-98-1	96
2			Ethanone, 1-phenyl-2-(4-pyridinyl)-	197	C13H11NO	001620-55-9	90
3			N-(5-Pyrimidinyl)benzamide	199	C11H9N3O	103854-65-5	62
4			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	62
5			Benzamide, N-(5-methyl-3-isoxazo...	202	C11H10N2O2	013053-85-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
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Instrument :
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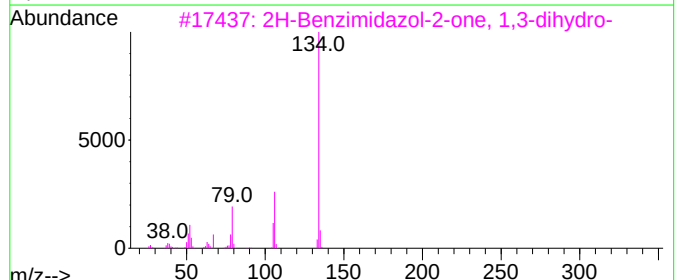
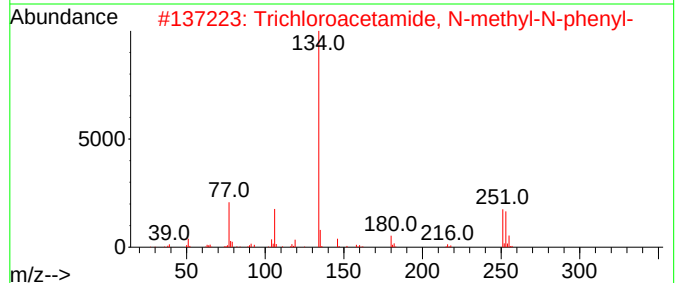
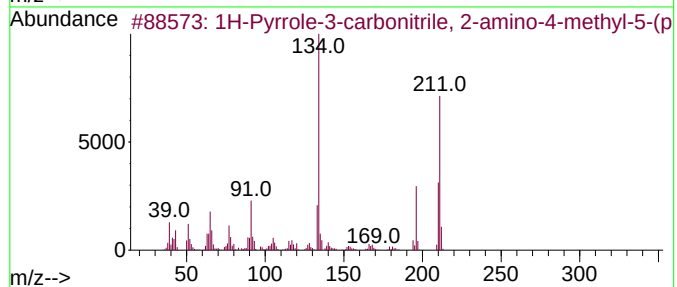
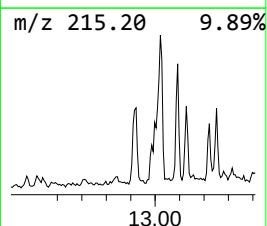
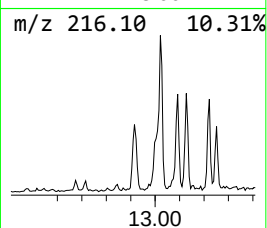
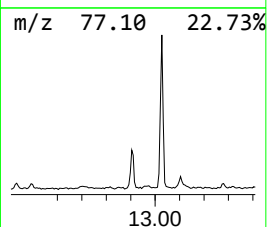
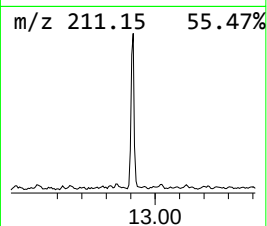
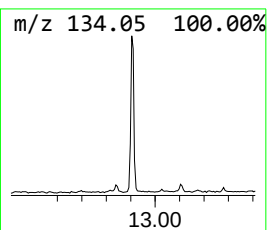
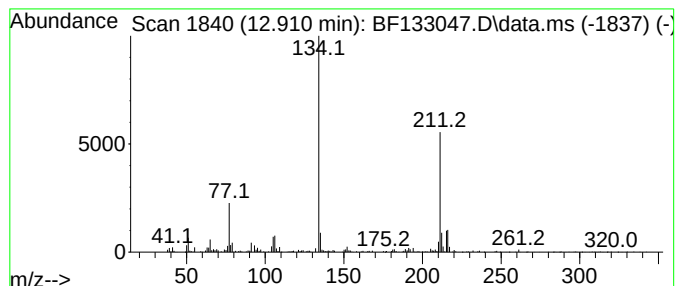
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Pyrrole-3-carbonitrile, ... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.910	3.67 ng	375122	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrrole-3-carbonitrile, 2-ami...	211	C13H13N3	1000351-31-6	53
2			Trichloroacetamide, N-methyl-N-p...	251	C9H8Cl3NO	1000345-68-4	37
3			2H-Benzimidazol-2-one, 1,3-dihydro-	134	C7H6N2O	000615-16-7	35
4			2-Benzoxazoline	134	C7H6N2O	004570-41-6	35
5			benzoic acid, 2-(methylamino)-, ...	272	C14H12N2O4	1000397-01-1	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
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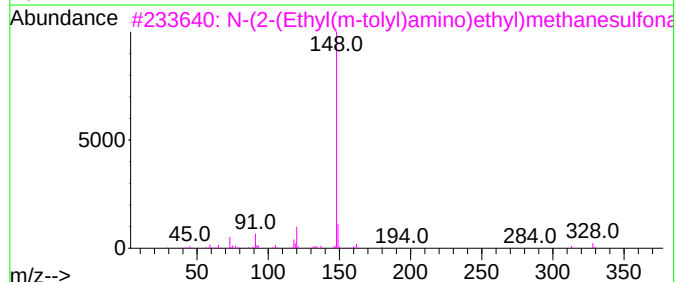
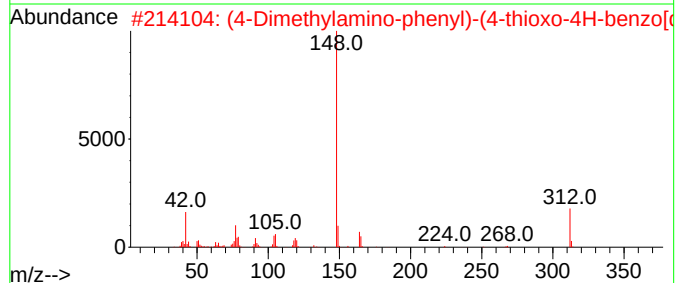
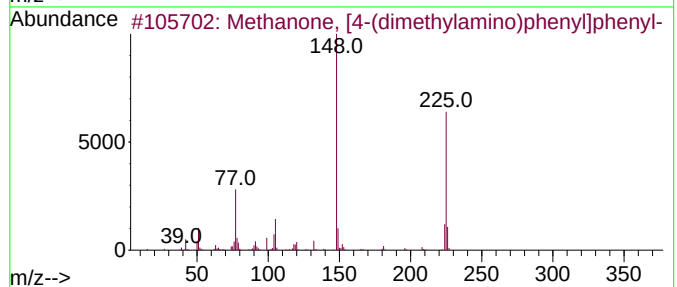
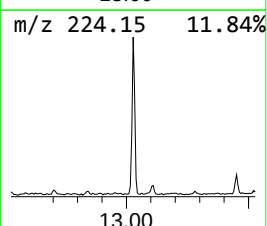
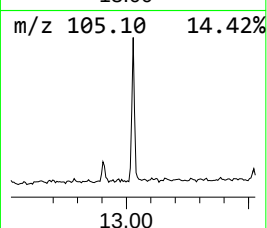
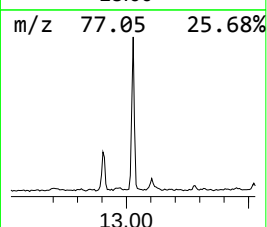
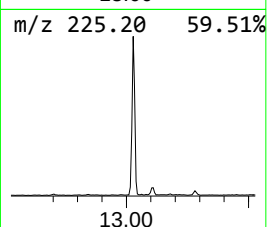
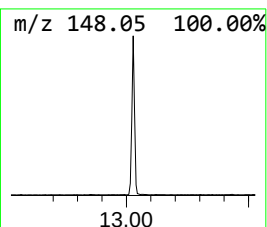
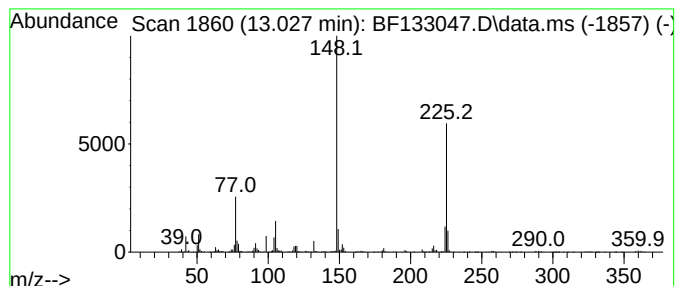
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Methanone, [4-(dimethylamin... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.027	13.98 ng	1429080	Chrysene-d12			13.892
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Methanone, [4-(dimethylamino)phe...		225	C15H15NO	000530-44-9	98
2	(4-Dimethylamino-phenyl)-(4-thio...		312	C17H16N2O2S	1000274-51-3	38
3	N-(2-(Ethyl(m-tolyl)amino)ethyl)...		328	C15H28N2O2SSi	1000505-33-6	38
4	3-Methylaniline, N-ethyl-N-pentyl-		205	C14H23N	1000310-30-2	38
5	4-(Dimethylamino)-N-(1,3-thiazol...		247	C12H13N3OS	851209-55-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
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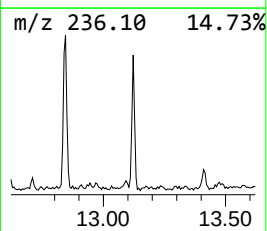
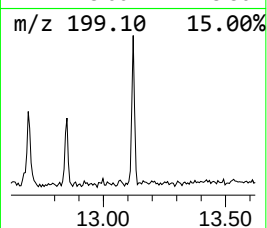
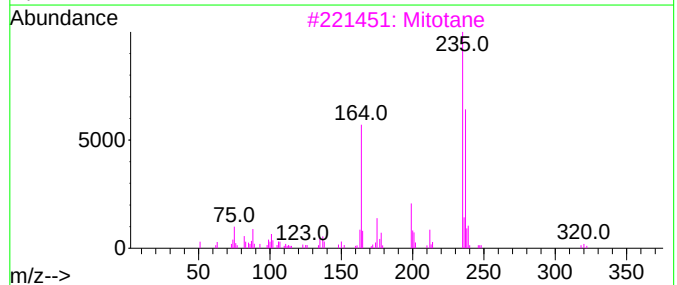
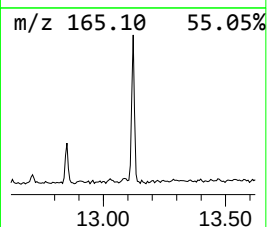
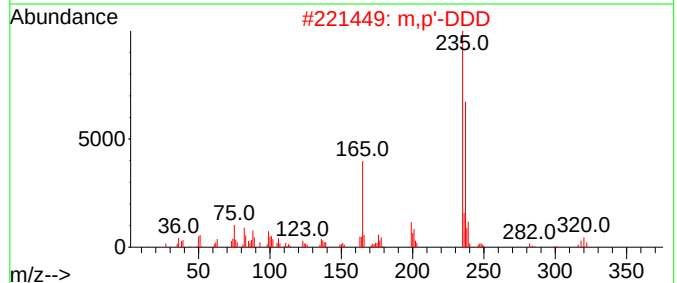
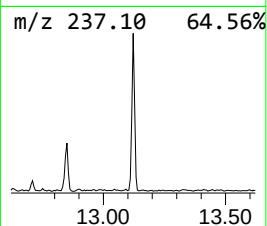
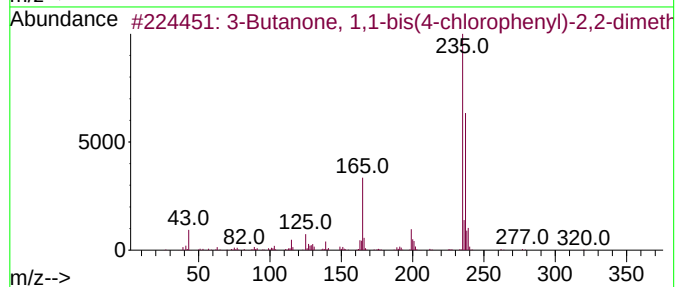
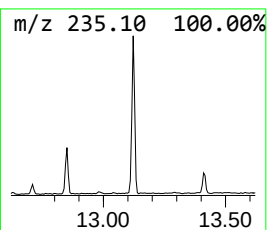
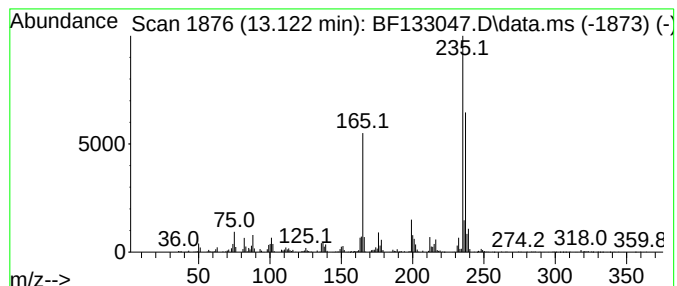
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 3-Butanone, 1,1-bis(4-chlor... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.122	6.57 ng	671489	Chrysene-d12	13.892

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	93
2		m,p'-DDD	318	C14H10Cl4	004329-12-8	93
3		Mitotane	318	C14H10Cl4	000053-19-0	89
4		2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	87
5		1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPC-3WC-23-04-17

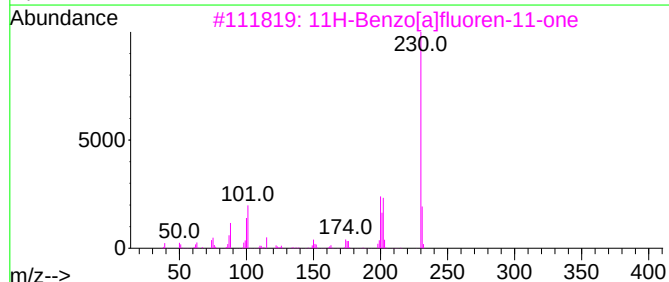
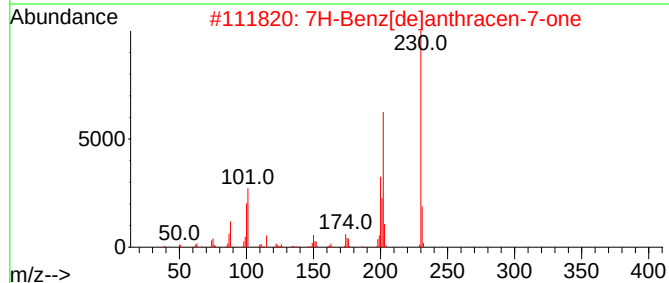
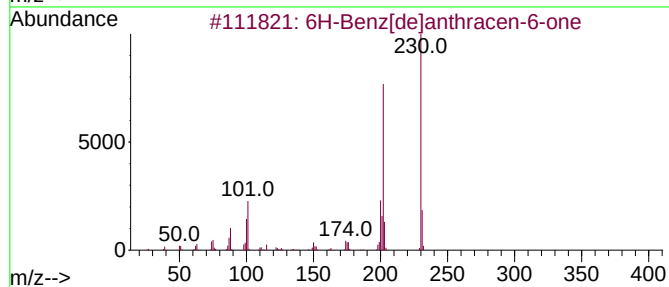
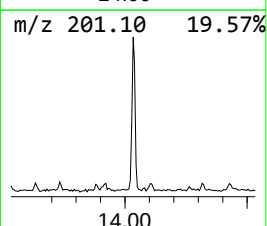
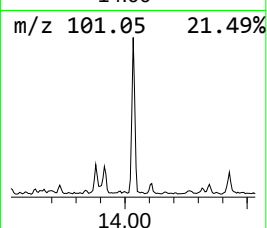
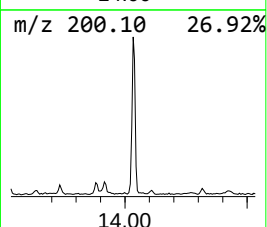
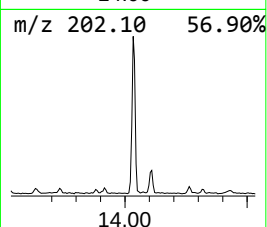
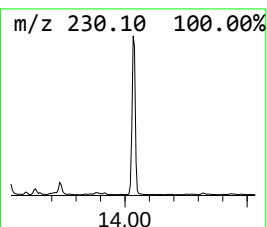
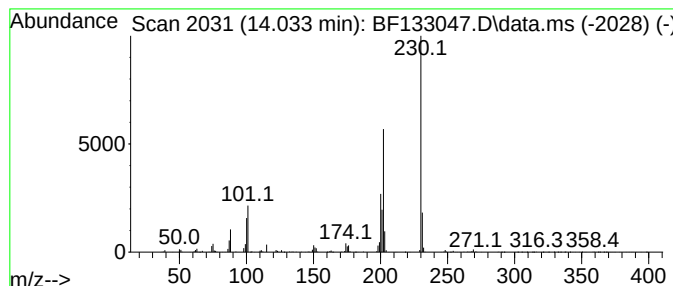
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 6H-Benz[de]anthracen-6-one Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.033	9.16 ng	936814	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	98
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	98
3			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
4			Pyrene	202	C16H10	000129-00-0	87
5			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPC-3WC-23-04-17

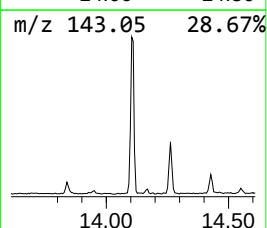
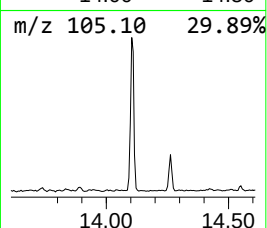
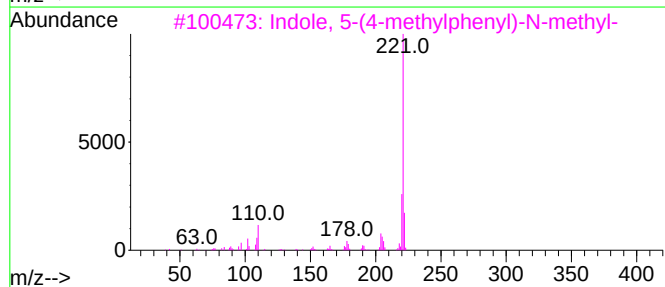
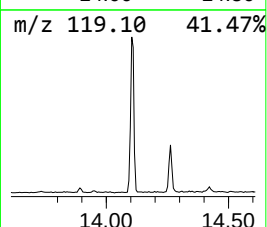
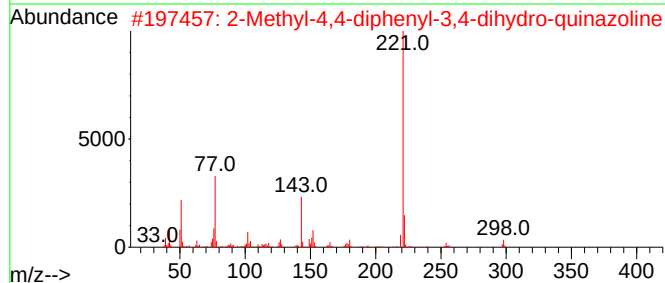
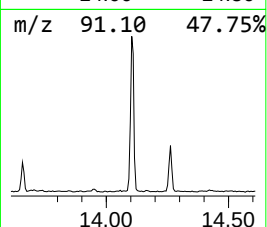
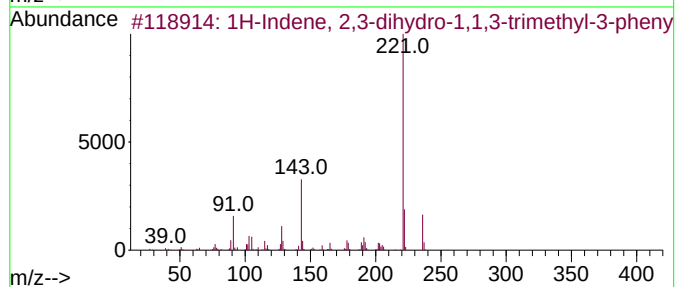
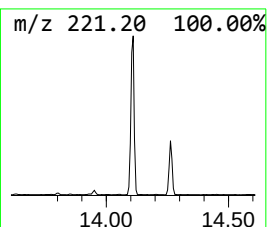
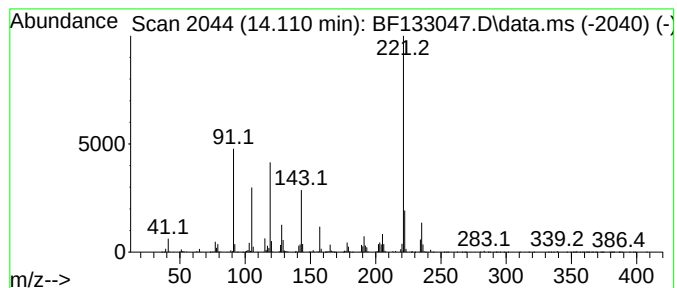
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown14.110 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.110	28.95 ng	2959110	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	47		
2	2-Methyl-4,4-diphenyl-3,4-dihydr...	298	C21H18N2	1000300-40-0	47		
3	Indole, 5-(4-methylphenyl)-N-met...	221	C16H15N	1000380-54-7	43		
4	1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	38		
5	2,3-Dihydro-2-methyl-4-phenyl-1H...	236	C16H16N2	111536-73-3	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SPC-3WC-23-04-17

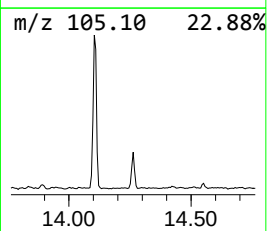
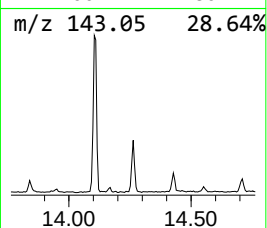
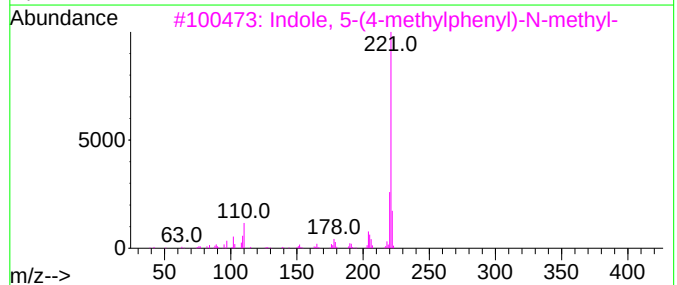
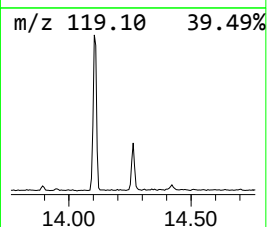
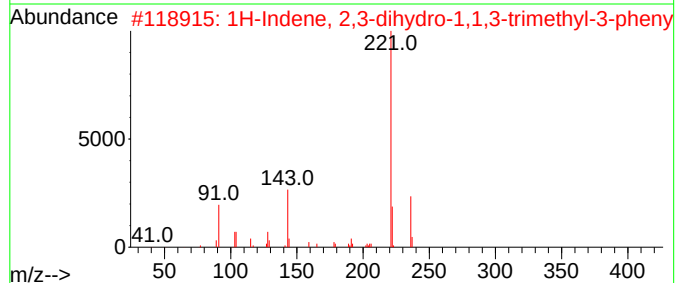
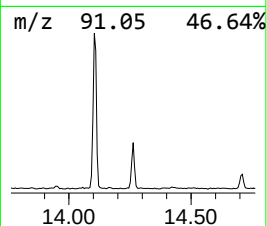
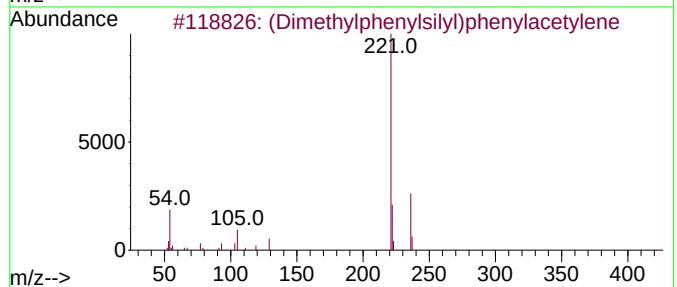
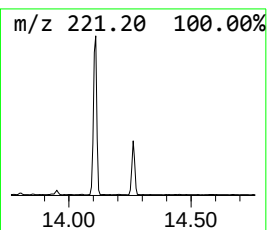
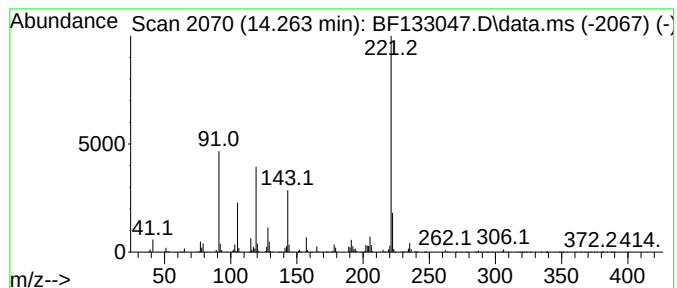
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown14.263 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.263	8.26 ng	844573	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(Dimethylphenylsilyl)phenylacety...	236	C16H16Si	1000434-31-1	47
2			1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	47
3			Indole, 5-(4-methylphenyl)-N-met...	221	C16H15N	1000380-54-7	38
4			1-(2-Hydroxy-4,5-dimethylphenyl)...	236	C13H20O2Si	1000484-73-7	35
5			1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

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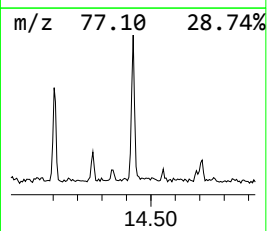
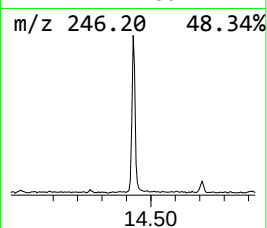
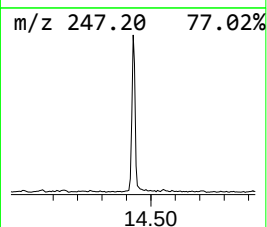
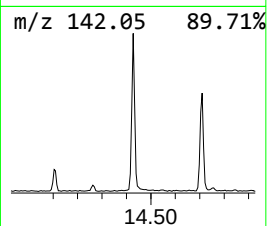
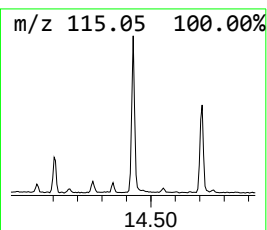
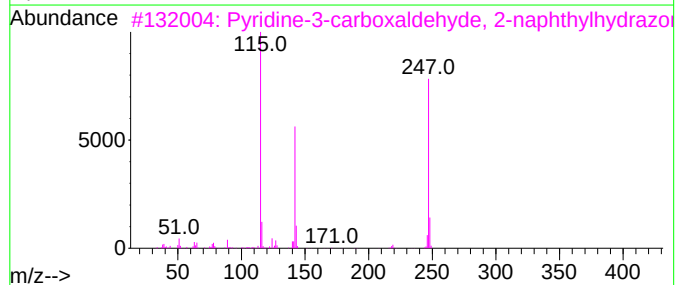
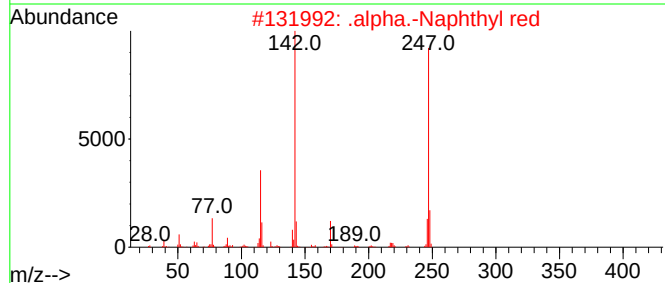
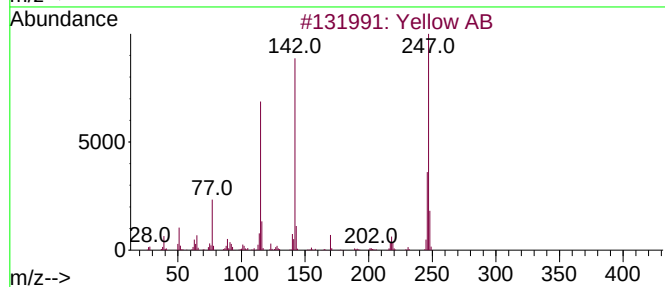
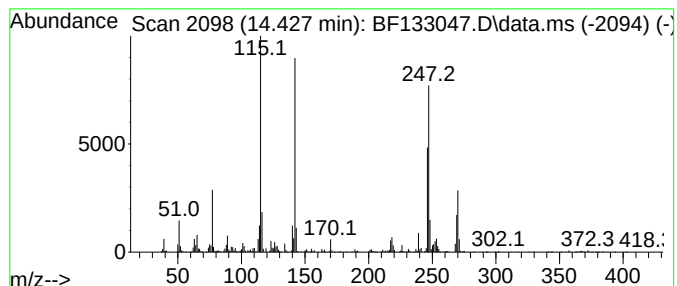
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Yellow AB Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.427	13.91 ng	1421470	Chrysene-d12	13.892

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Yellow AB		247	C16H13N3	000085-84-7	93
2	.alpha.-Naphthyl red		247	C16H13N3	000131-22-6	60
3	Pyridine-3-carboxaldehyde, 2-nap...		247	C16H13N3	1000262-93-3	53
4	Naphthalene, 1-methyl-		142	C11H10	000090-12-0	50
5	3-Methyl-3H-naphth[1,2-e]indol-1...		247	C17H13NO	098033-21-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SPC-3WC-23-04-17

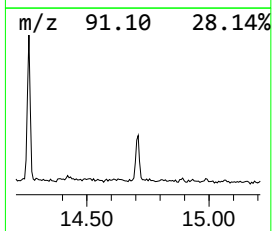
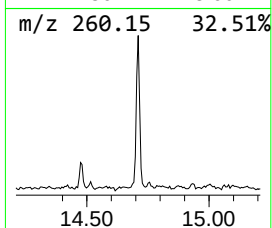
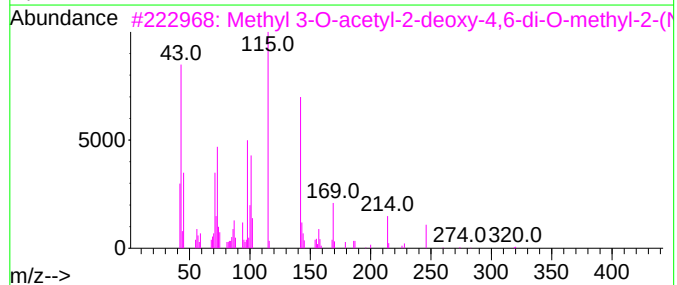
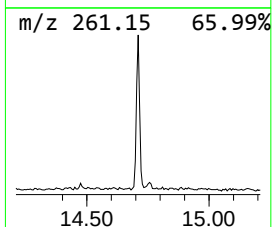
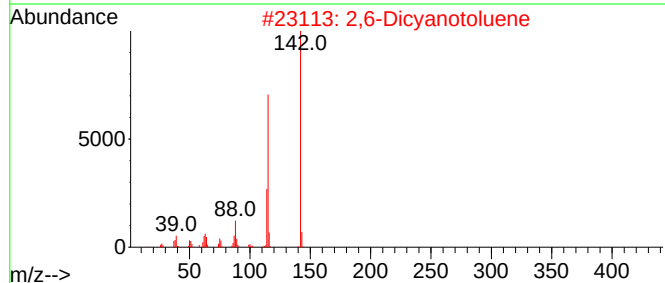
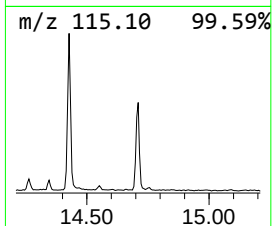
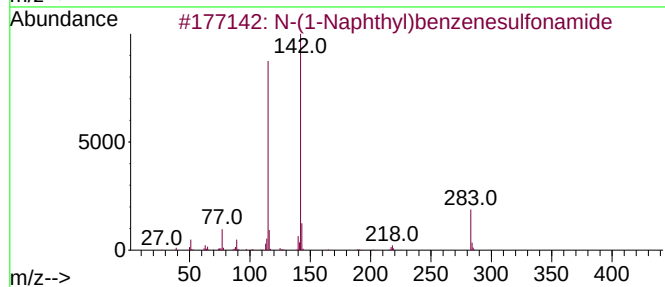
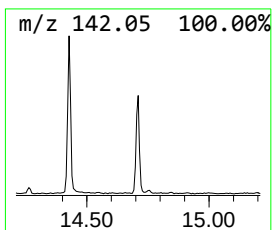
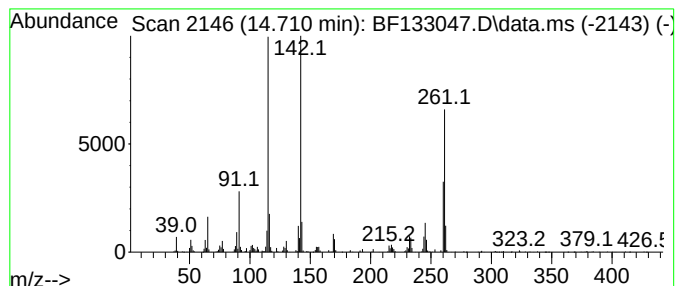
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 N-(1-Naphthyl)benzenesulfon... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.710	13.87 ng	653045	Perylene-d12	15.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-(1-Naphthyl)benzenesulfonamide	283	C16H13NO2S	015309-82-7	50
2			2,6-Dicyanotoluene	142	C9H6N2	002317-22-8	49
3			Methyl 3-O-acetyl-2-deoxy-4,6-di...	319	C14H25NO7	056389-83-4	47
4			Benzamide, 2-nitro-N-(1-naphthal...	292	C17H12N2O3	038267-68-4	40
5			5-Methylisophthalonitrile	142	C9H6N2	039718-07-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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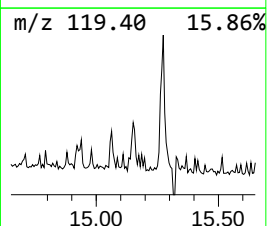
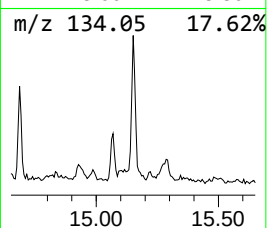
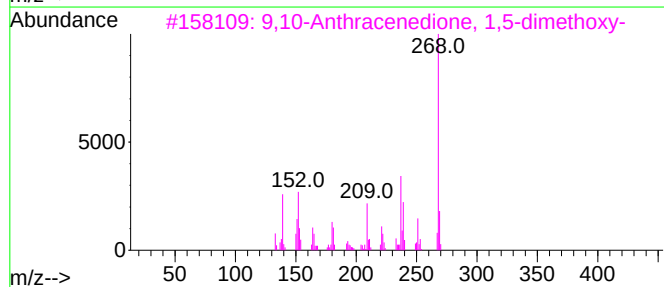
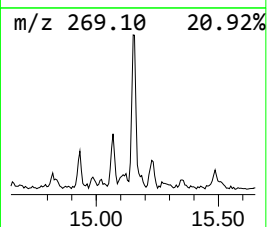
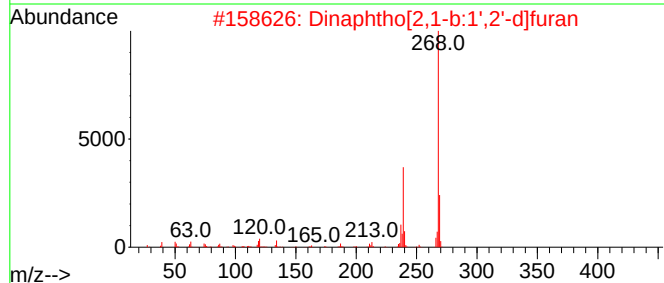
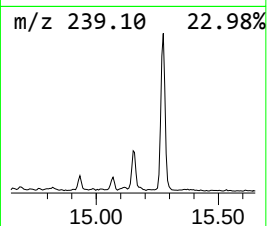
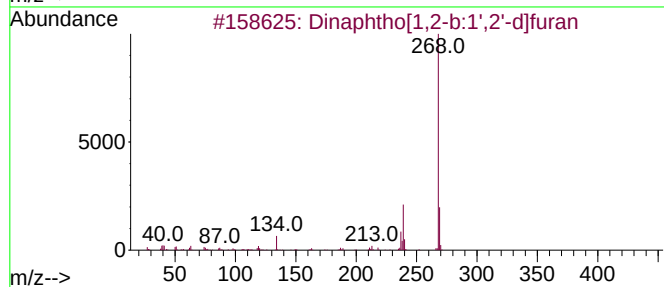
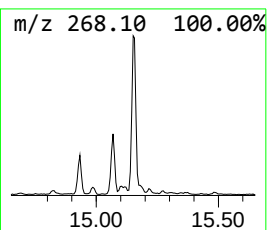
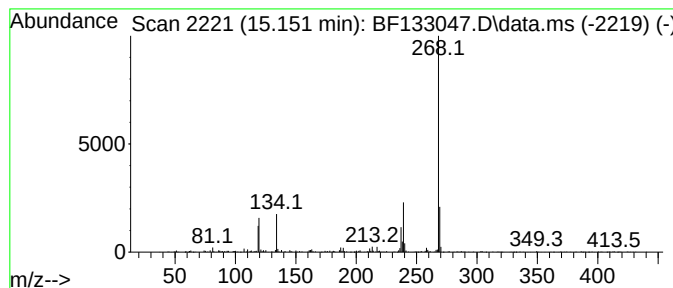
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Dinaphtho[1,2-b:1',2'-d]furan Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.151	9.72 ng	457524	Perylene-d12	15.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dinaphtho[1,2-b:1',2'-d]furan	268	C20H12O	000207-93-2	91
2			Dinaphtho[2,1-b:1',2'-d]furan	268	C20H12O	000194-63-8	87
3			9,10-Anthracenedione, 1,5-dimeth...	268	C16H12O4	006448-90-4	64
4			Benzo(a)pyrene 4,5-oxide	268	C20H12O	037574-47-3	64
5			7-Hydroxy-3-(2-methoxyphenyl)cou...	268	C16H12O4	549501-06-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPC-3WC-23-04-17

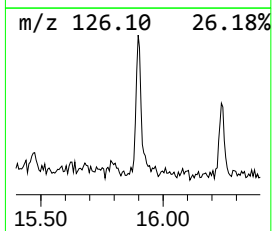
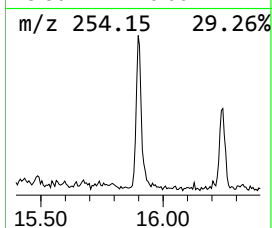
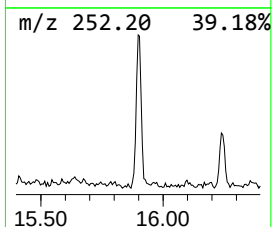
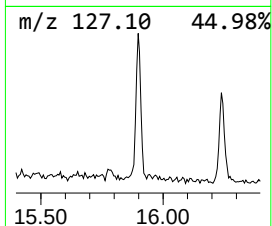
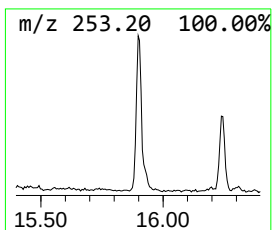
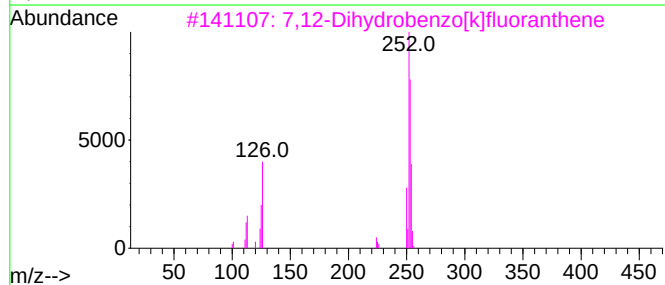
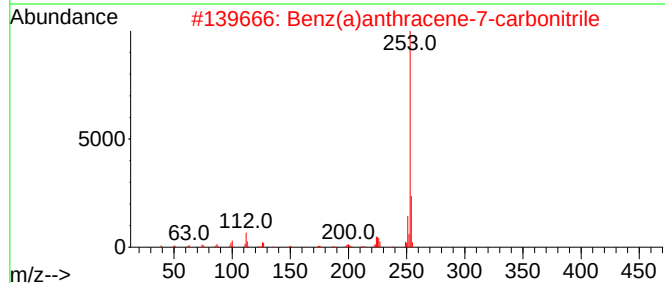
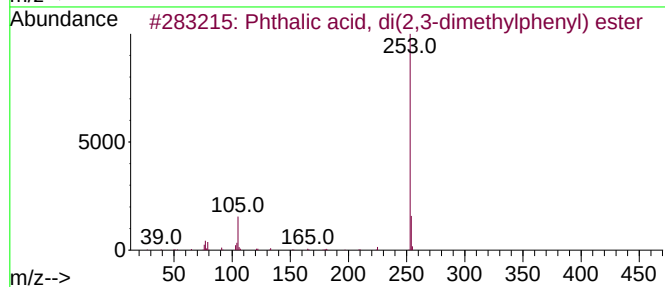
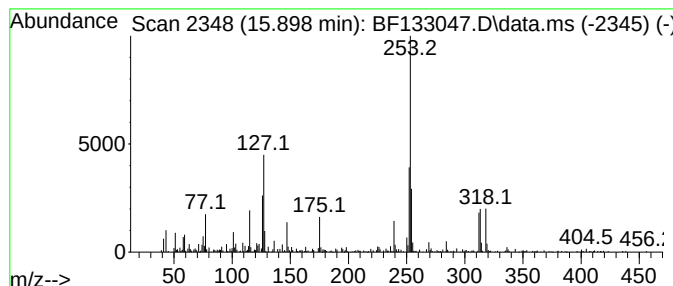
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown15.898 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.898	6.74 ng	317518	Perylene-d12	15.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, di(2,3-dimethylph...	374	C24H22O4	1000357-09-2	22
2			Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	16
3			7,12-Dihydrobenzo[k]fluoranthene	254	C20H14	1000080-17-7	14
4			1H-Pyrrolo[2,3-f]quinolin-9-ol, ...	254	C16H18N2O	1000303-71-9	14
5			3-Chloro-4-iodoaniline	253	C6H5ClIN	135050-44-1	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPC-3WC-23-04-17

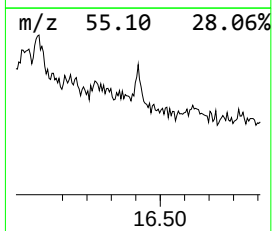
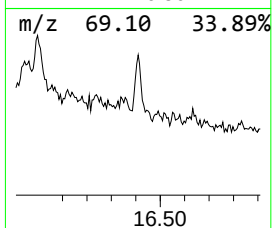
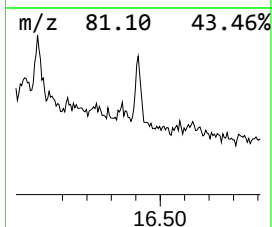
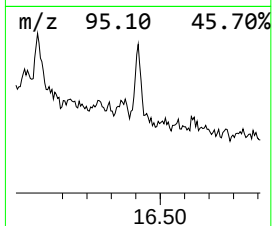
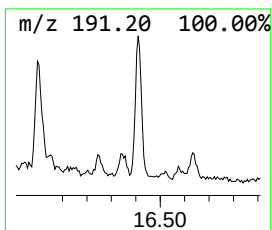
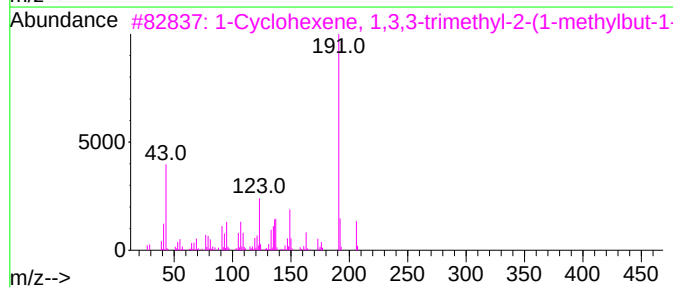
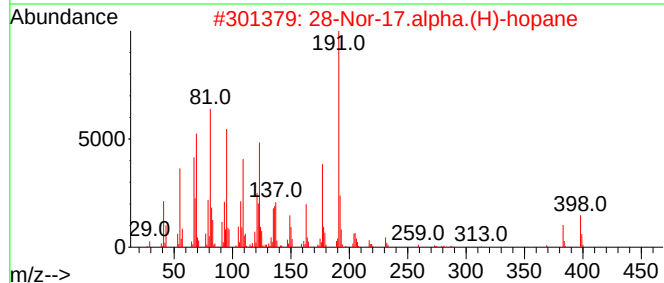
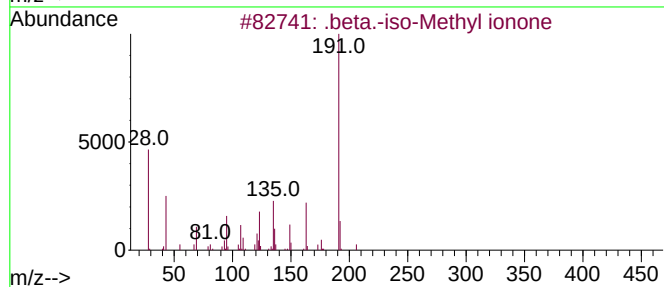
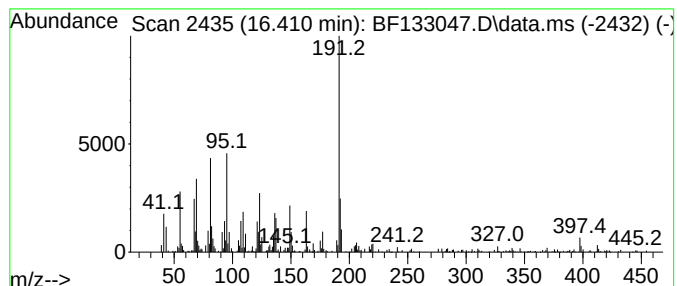
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 .beta.-iso-Methyl ionone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.410	9.56 ng	450096	Perylene-d12	15.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	74
2		28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	53
3		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	53
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	49
5		17.alpha.(H),21.beta.(H)-Hopane	412	C30H52	013849-96-2	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

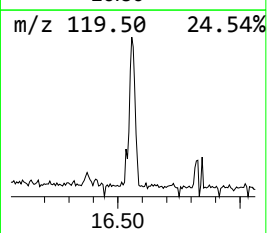
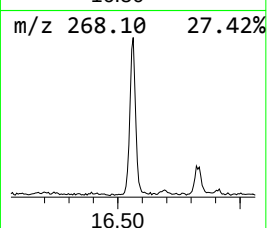
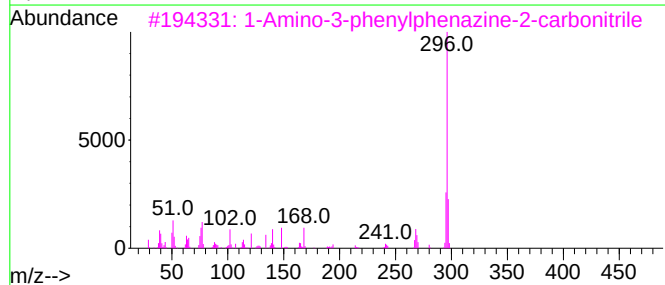
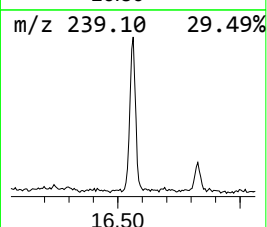
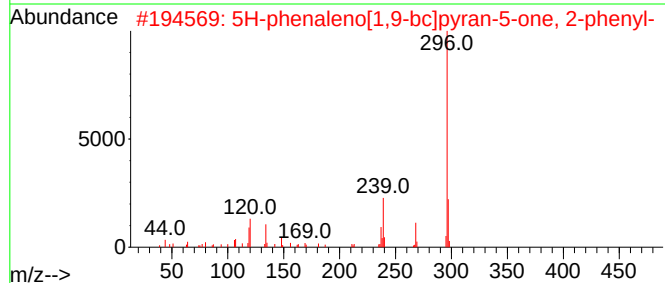
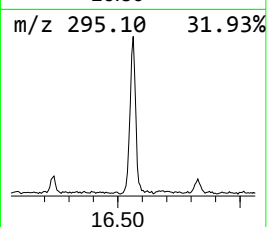
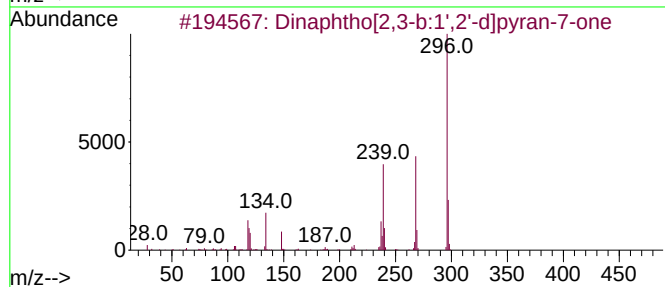
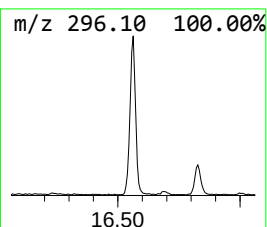
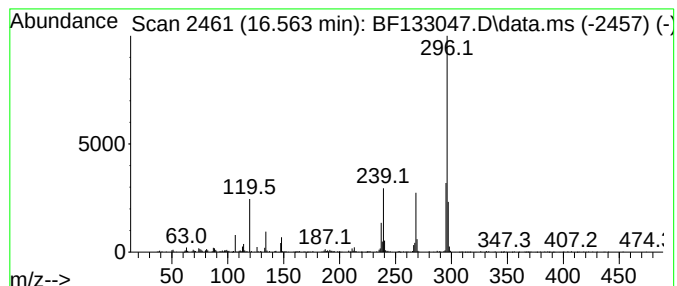
Instrument :
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ClientSampleId :
SPC-3WC-23-04-17

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Dinaphtho[2,3-b:1',2'-d]pyr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.563	12.63 ng	594550	Perylene-d12		15.321	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dinaphtho[2,3-b:1',2'-d]pyran-7-one		296	C21H12O2	154352-87-1	95
2	5H-phenaleno[1,9-bc]pyran-5-one,...		296	C21H12O2	1000400-06-2	81
3	1-Amino-3-phenylphenazine-2-carb...		296	C19H12N4	1000444-09-5	58
4	Imidazolidine-2,4-dione, 5-(4-et...		296	C17H16N2O3	086029-57-4	58
5	7-Ethoxy-3-(4-methoxyphenyl)coum...		296	C18H16O4	263364-70-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPC-3WC-23-04-17

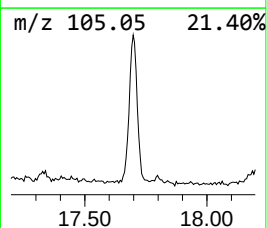
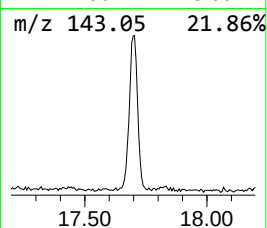
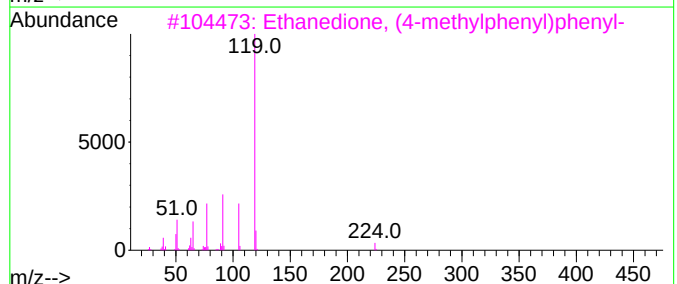
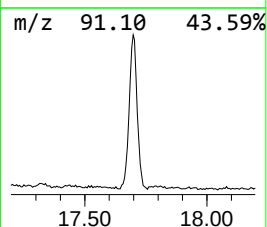
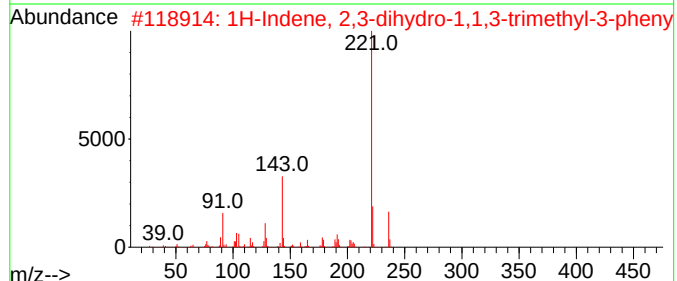
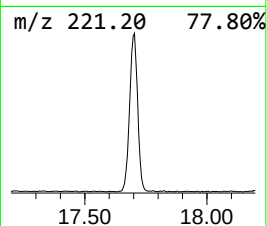
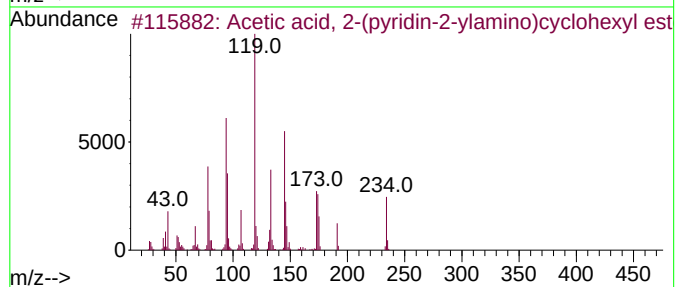
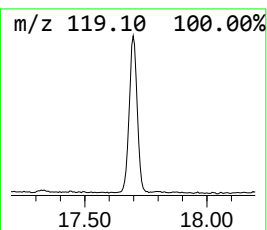
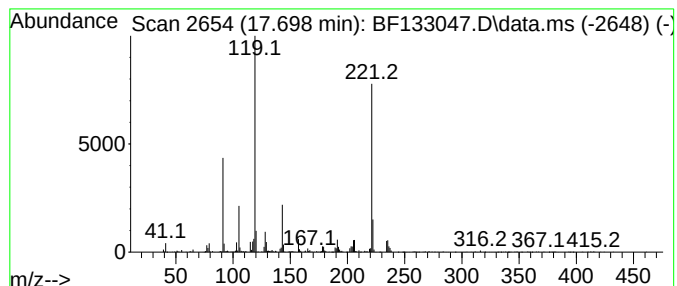
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Acetic acid, 2-(pyridin-2-y... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.698	29.31 ng	1379810	Perylene-d12	15.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 2-(pyridin-2-ylamin...	234	C13H18N2O2	1000185-40-1	83
2			1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	38
3			Ethanedione, (4-methylphenyl)phe...	224	C15H12O2	002431-00-7	35
4			.alpha.,.alpha.-DIMETHYLBENZYL I...	206	C13H18O2	1000429-51-9	25
5			Indole, 5-(4-methylphenyl)-N-met...	221	C16H15N	1000380-54-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133047.D
Acq On : 26 Apr 2023 01:34
Operator : CG\JU
Sample : 02390-11 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.422	7.422	5.4	ng	455177	2	8.022	1678130	20.0
2-Aminobiphenyl	10.286	6.3	ng	589067	3	9.775	1865300	20.0
1H-Indene, 2,3-...	10.869	3.5	ng	291243	4	11.257	1659510	20.0
7-Acetyl-6-ethy...	11.186	4.0	ng	333120	4	11.257	1659510	20.0
n-Hexadecanoic ...	11.792	8.7	ng	725609	4	11.257	1659510	20.0
2,2',4,5-Tetrac...	11.875	3.9	ng	324196	4	11.257	1659510	20.0
Benzamide, N-ph...	11.928	10.7	ng	883916	4	11.257	1659510	20.0
1H-Pyrrole-3-ca...	12.910	3.7	ng	375122	5	13.892	2044430	20.0
Methanone, [4-(...	13.027	14.0	ng	1429080	5	13.892	2044430	20.0
3-Butanone, 1,1...	13.122	6.6	ng	671489	5	13.892	2044430	20.0
6H-Benz[de]anth...	14.033	9.2	ng	936814	5	13.892	2044430	20.0
unknown14.110	14.110	28.9	ng	2959110	5	13.892	2044430	20.0
unknown14.263	14.263	8.3	ng	844573	5	13.892	2044430	20.0
Yellow AB	14.427	13.9	ng	1421470	5	13.892	2044430	20.0
N-(1-Naphthyl)b...	14.710	13.9	ng	653045	6	15.321	941622	20.0
Dinaphtho[1,2-b...	15.151	9.7	ng	457524	6	15.321	941622	20.0
unknown15.898	15.898	6.7	ng	317518	6	15.321	941622	20.0
.beta.-iso-Meth...	16.410	9.6	ng	450096	6	15.321	941622	20.0
Dinaphtho[2,3-b...	16.563	12.6	ng	594550	6	15.321	941622	20.0
Acetic acid, 2-...	17.698	29.3	ng	1379810	6	15.321	941622	20.0