

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
 Data File : BF086110.D
 Acq On : 6 Apr 2016 21:01
 Operator : UM/SJ
 Sample : H2074-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB01A

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-BF040216.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	2.201	6	10	11	rBV4	151412	323543	11.82%	0.677%
2	2.350	21	23	26	rVB3	145032	216742	7.92%	0.453%
3	2.407	26	28	30	rBV3	106479	219277	8.01%	0.459%
4	2.476	33	34	35	rBV	218041	204396	7.47%	0.428%
5	2.510	35	37	40	rBV4	61475	107815	3.94%	0.226%
6	2.613	45	46	48	rVB	334593	239020	8.74%	0.500%
7	2.727	52	56	58	rBV3	349682	1160104	42.40%	2.427%
8	2.761	58	59	60	rVV	139278	136845	5.00%	0.286%
9	2.784	60	61	63	rVB2	561269	438299	16.02%	0.917%
10	2.887	69	70	72	rBV	336675	652451	23.84%	1.365%
11	2.978	77	78	79	rVV	154064	191922	7.01%	0.402%
12	3.070	82	86	87	rVV4	162538	397375	14.52%	0.831%
13	3.093	87	88	90	rVV2	436949	418883	15.31%	0.876%
14	3.207	97	98	100	rBV2	70341	97654	3.57%	0.204%
15	3.264	102	103	105	rVB2	224992	207038	7.57%	0.433%
16	3.298	105	106	108	rBV2	141521	268349	9.81%	0.561%
17	3.378	111	113	115	rBV3	69700	117415	4.29%	0.246%
18	3.573	127	130	132	rBV4	73111	157818	5.77%	0.330%
19	3.744	141	145	146	rBV3	194127	536862	19.62%	1.123%
20	3.779	146	148	152	rVV5	290463	776348	28.37%	1.624%
21	3.858	152	155	158	rVV5	229414	755077	27.60%	1.580%
22	3.916	158	160	161	rVV2	205650	412636	15.08%	0.863%
23	3.996	164	167	171	rVV4	1276535	2736274	100.00%	5.725%
24	4.076	173	174	177	rVV3	323366	310774	11.36%	0.650%
25	4.133	177	179	182	rVB3	350119	546148	19.96%	1.143%
26	4.201	182	185	186	rBV3	355893	671872	24.55%	1.406%
27	4.224	186	187	189	rVB	402107	290012	10.60%	0.607%
28	4.910	245	247	253	rBV	73232	123486	4.51%	0.258%
29	5.344	282	285	287	rBV	1600759	2286020	83.54%	4.783%
30	6.396	375	377	379	rBV	2154407	2387688	87.26%	4.995%
31	6.510	385	387	390	rBV	2192415	2259952	82.59%	4.728%
32	6.739	405	407	411	rVB	708220	704793	25.76%	1.474%
33	6.899	419	421	423	rBV	2155612	1963533	71.76%	4.108%
34	7.310	454	457	460	rBV	1532786	1426040	52.12%	2.983%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	8.030	517	520	522 rBV	929557	868773	31.75%	1.818%
36	9.105	611	614	617 rBV	2668431	2372112	86.69%	4.963%
37	9.505	647	649	651 rBV	385373	336834	12.31%	0.705%
38	9.779	671	673	675 rVV	934898	833404	30.46%	1.744%
39	10.213	710	711	713 rVB	104798	76898	2.81%	0.161%
40	10.328	719	721	724 rVB	100109	99364	3.63%	0.208%
41	10.431	724	730	733 rBV3	46438	141975	5.19%	0.297%
42	10.568	739	742	745 rBV	1294018	1253819	45.82%	2.623%
43	10.682	751	752	758 rVB	99674	179173	6.55%	0.375%
44	10.876	766	769	770 rVV2	54968	81199	2.97%	0.170%
45	11.002	778	780	784 rVV2	40867	103043	3.77%	0.216%
46	11.071	784	786	788 rVV	43921	81702	2.99%	0.171%
47	11.128	789	791	793 rVV2	77907	145190	5.31%	0.304%
48	11.162	793	794	797 rVB	99901	89046	3.25%	0.186%
49	11.265	801	803	804 rBV	682199	845231	30.89%	1.768%
50	11.288	804	805	807 rVV	973911	713597	26.08%	1.493%
51	11.334	807	809	811 rVV	195445	242867	8.88%	0.508%
52	11.505	821	824	827 rVV2	62986	129780	4.74%	0.272%
53	11.562	827	829	830 rVV	68204	92507	3.38%	0.194%
54	11.585	830	831	836 rVV2	38171	108844	3.98%	0.228%
55	11.768	844	847	848 rVV	131445	205987	7.53%	0.431%
56	11.791	848	849	852 rVV2	281464	373071	13.63%	0.780%
57	11.871	855	856	861 rVV	208544	320546	11.71%	0.671%
58	11.962	861	864	866 rVB2	39150	77947	2.85%	0.163%
59	12.054	868	872	874 rBV4	77399	110159	4.03%	0.230%
60	12.202	883	885	887 rBV3	57619	102003	3.73%	0.213%
61	12.236	887	888	893 rVV	69122	154868	5.66%	0.324%
62	12.316	893	895	896 rVV	97374	132779	4.85%	0.278%
63	12.351	896	898	901 rVV	113836	172948	6.32%	0.362%
64	12.408	901	903	906 rVV	75595	117098	4.28%	0.245%
65	12.477	906	909	911 rVV	1326790	1269270	46.39%	2.655%
66	12.534	911	914	916 rVV3	64675	131746	4.81%	0.276%
67	12.579	916	918	920 rVV2	137231	177131	6.47%	0.371%
68	12.625	920	922	923 rVV2	73108	98345	3.59%	0.206%
69	12.705	926	929	931 rBV	1014350	1347934	49.26%	2.820%
70	12.842	939	941	944 rVV	1606372	1736143	63.45%	3.632%
71	12.922	945	948	954 rVB3	92085	209118	7.64%	0.437%

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72	13.025	954	957	959	rBV2	137599	296355	10.83%	0.620%
73	13.094	959	963	965	rBV2	97205	166516	6.09%	0.348%
74	13.128	965	966	970	rVB	94364	146451	5.35%	0.306%
75	13.254	976	977	981	rBV	82628	105041	3.84%	0.220%
76	13.425	986	992	993	rBV4	78951	200248	7.32%	0.419%
77	13.539	1001	1002	1004	rVB	97239	98677	3.61%	0.206%
78	13.654	1009	1012	1013	rBV	99237	184243	6.73%	0.385%
79	13.699	1015	1016	1019	rVV2	159498	182551	6.67%	0.382%
80	13.745	1019	1020	1023	rVV	266605	336138	12.28%	0.703%
81	13.894	1030	1033	1035	rBV2	896392	1528271	55.85%	3.197%
82	13.997	1040	1042	1046	rVB2	74752	143447	5.24%	0.300%
83	14.065	1046	1048	1050	rBV3	86378	154827	5.66%	0.324%
84	14.282	1065	1067	1069	rBV	90573	127922	4.68%	0.268%
85	14.362	1072	1074	1077	rBV3	54254	131380	4.80%	0.275%
86	14.602	1092	1095	1098	rBV3	112884	293739	10.73%	0.615%
87	14.911	1119	1122	1126	rBV	550346	1000093	36.55%	2.092%
88	14.968	1126	1127	1129	rVV2	84912	92195	3.37%	0.193%
89	15.014	1129	1131	1136	rVB	93841	171937	6.28%	0.360%
90	15.185	1144	1146	1149	rVB	301826	337275	12.33%	0.706%
91	15.243	1149	1151	1154	rVV	408423	517667	18.92%	1.083%
92	15.300	1154	1156	1162	rVB2	468462	749647	27.40%	1.568%
93	15.848	1202	1204	1207	rVB4	70350	109087	3.99%	0.228%
94	16.294	1241	1243	1251	rVB5	63375	160236	5.86%	0.335%
95	16.648	1271	1274	1278	rVB	146463	297691	10.88%	0.623%
96	16.820	1286	1289	1291	rBV2	61775	154684	5.65%	0.324%
97	17.048	1306	1309	1314	rVB	100029	200313	7.32%	0.419%
98	17.780	1371	1373	1376	rBV4	36052	91757	3.35%	0.192%
99	18.260	1412	1415	1418	rVB4	176099	356965	13.05%	0.747%
100	18.614	1443	1446	1452	rBV7	46586	188745	6.90%	0.395%

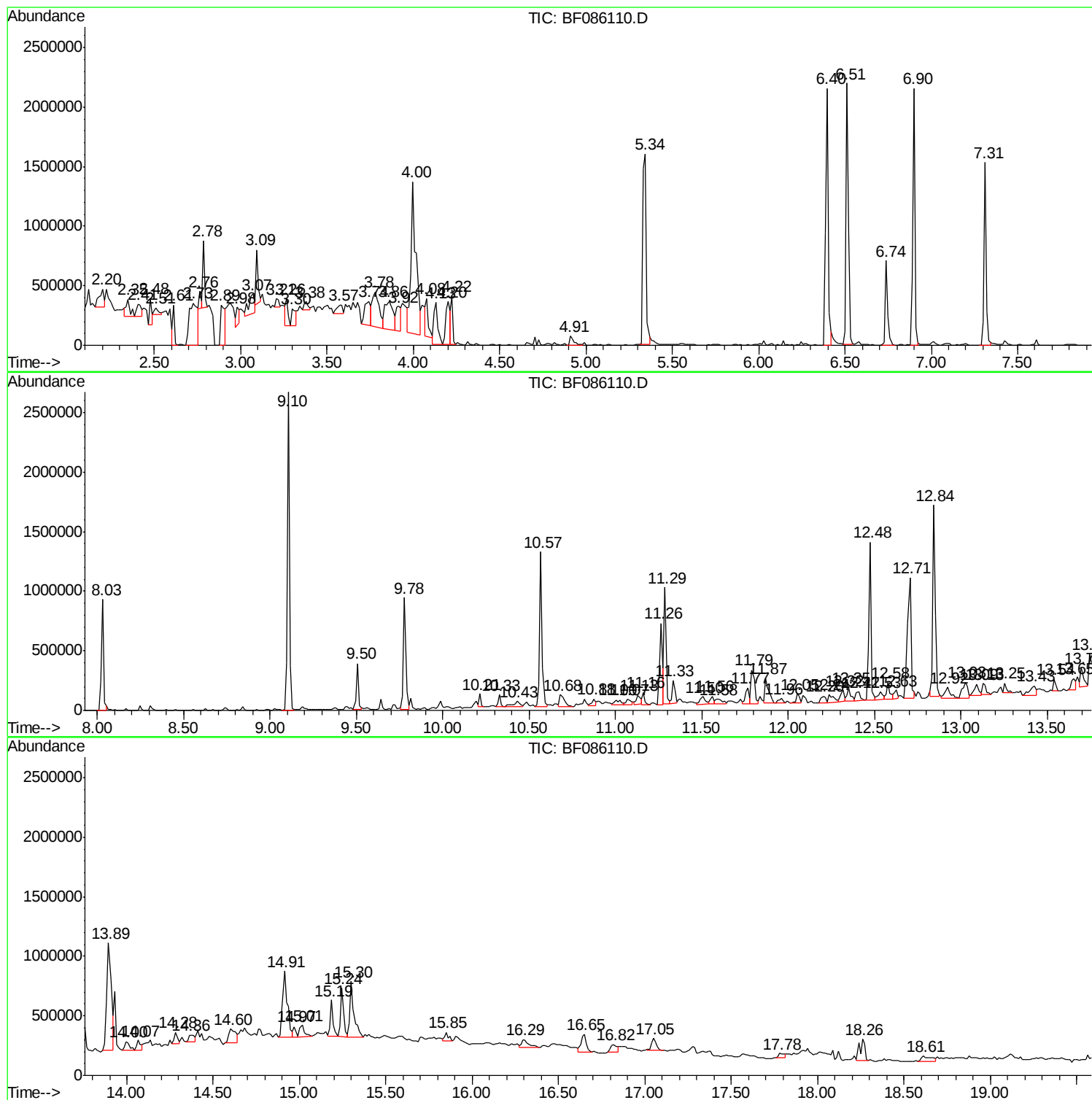
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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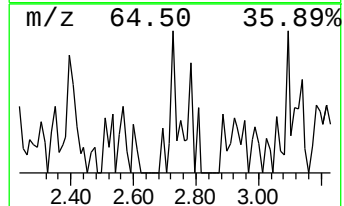
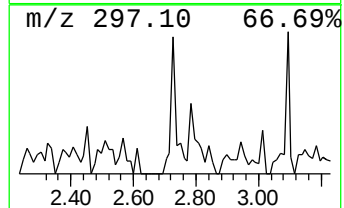
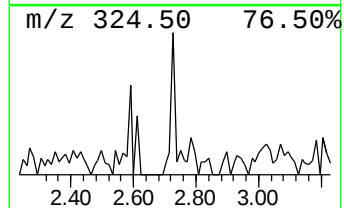
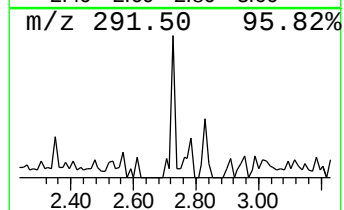
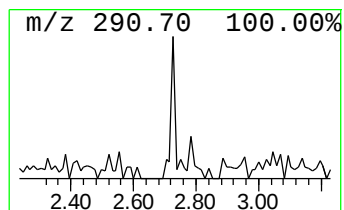
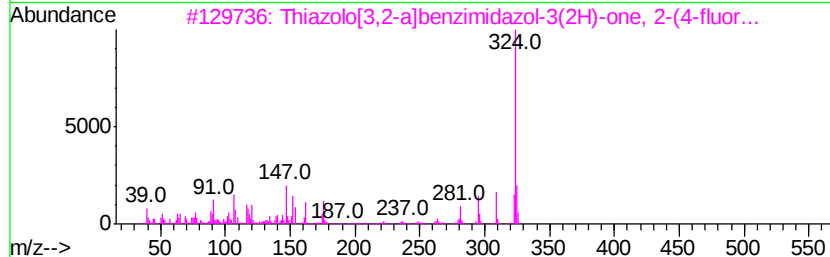
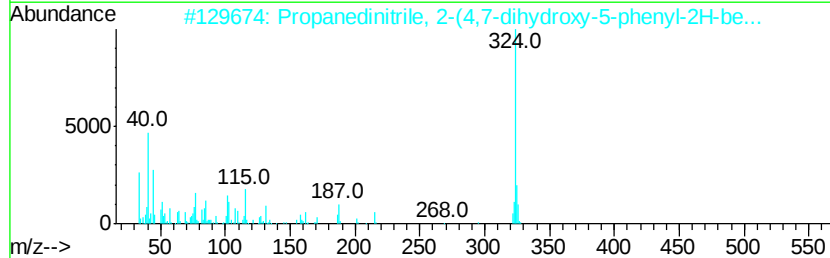
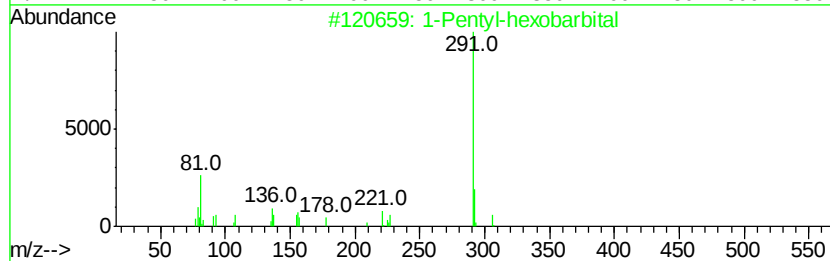
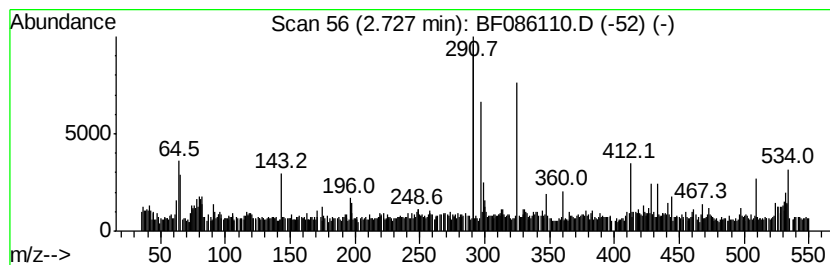
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown2.73 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.73	32.92 ng	1160100	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentyl-hexobarbital	306	C17H26N2O3	1000100-94-9	25
2		Propanedinitrile, 2-(4,7-dihydro...	324	C16H8N2O2S2	1000269-50-5	25
3		Thiazolo[3,2-a]benzimidazol-3(2H)...	324	C18H13FN2OS	294878-43-6	25
4		Thiazolo[3,2-a]benzimidazol-3(2H)...	324	C18H13FN2OS	294878-41-4	25
5		3-Fluoro-3-(4-hydroxy-3,5-dimeth...	324	C13H12F4O3S	1000276-10-2	20



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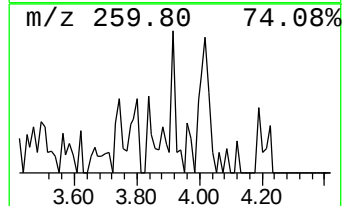
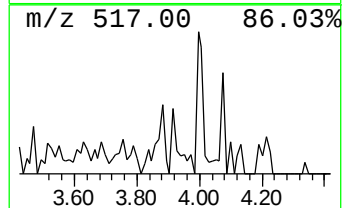
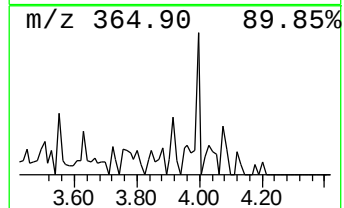
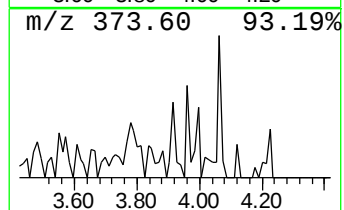
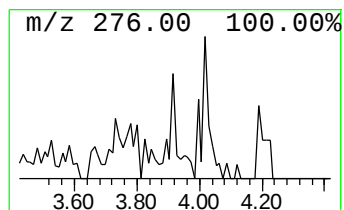
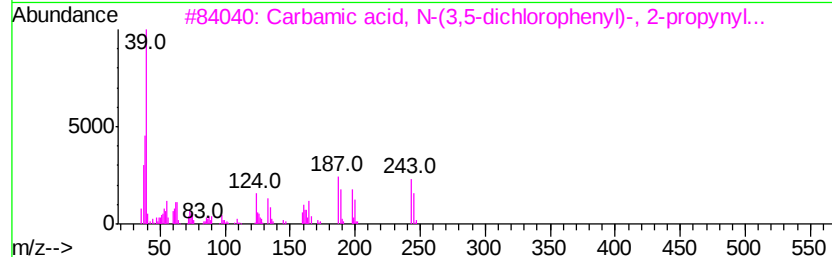
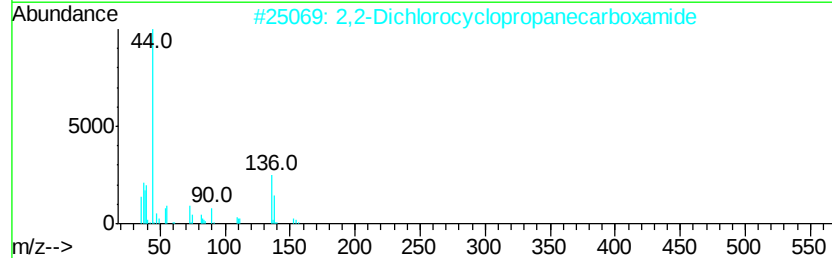
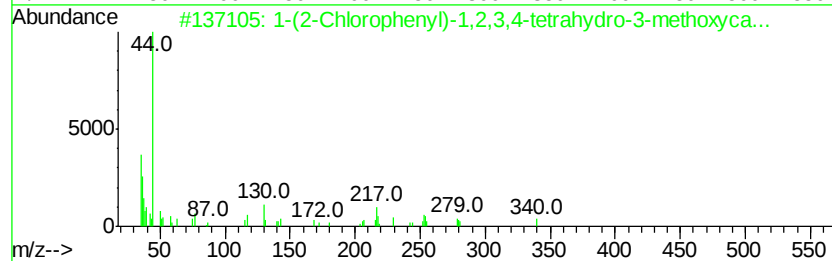
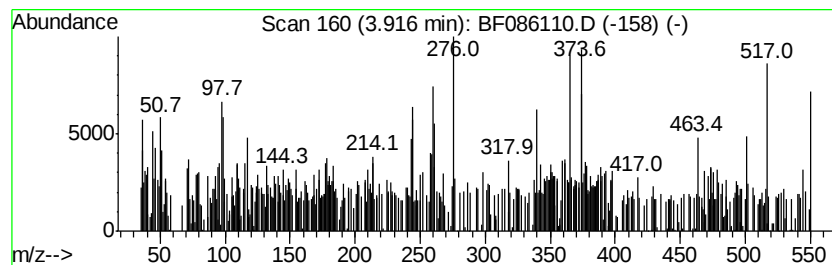
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown3.92 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	11.71 ng	412636	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(2-Chlorophenyl)-1,2,3,4-tetra...	340	C19H17ClN2O2	223690-84-4	23
2		2,2-Dichlorocyclopropanecarboxamide	153	C4H5Cl2NO	075885-60-8	9
3		Carbamic acid, N-(3,5-dichloroph...	243	C10H7Cl2NO2	025216-52-8	7
4		1,3,5,7,9,11,13-Heptaethyl-5,9,1...	546	C14H38O9Si7	080177-43-1	7
5		Beryllium, hexakis(chloroacetato...	610	C12H12Be4Cl6O13	012560-98-4	7



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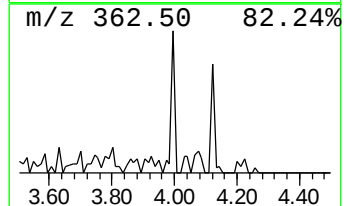
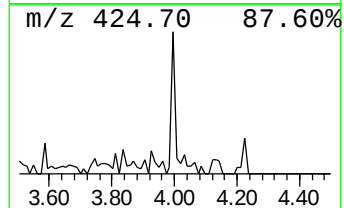
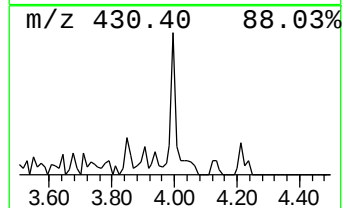
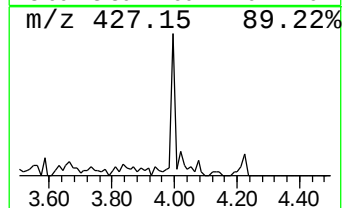
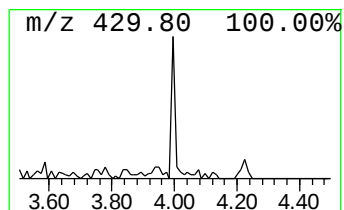
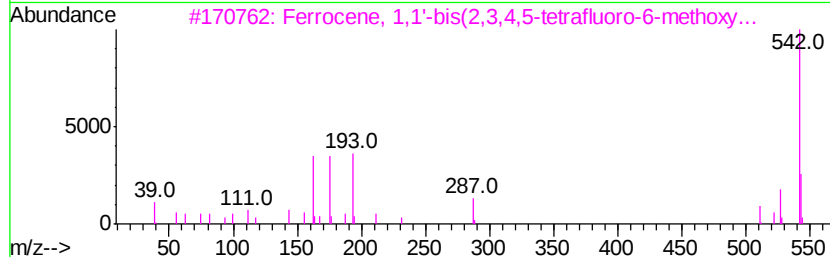
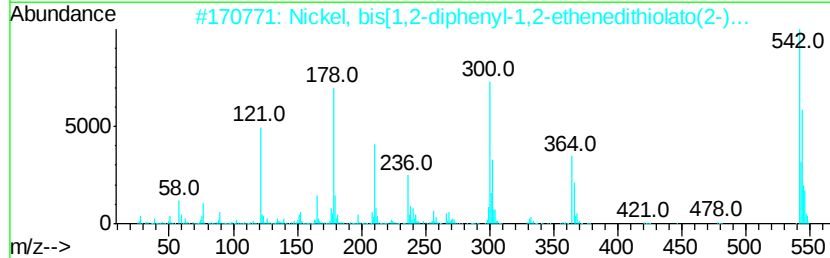
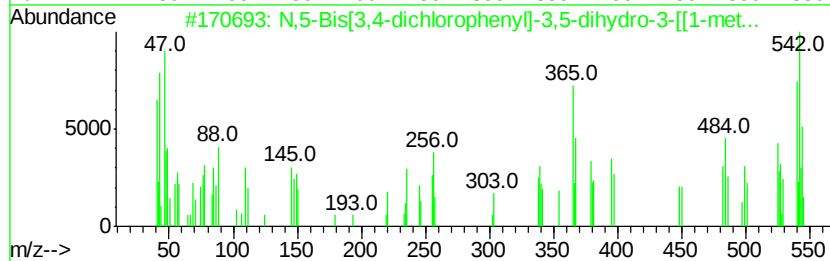
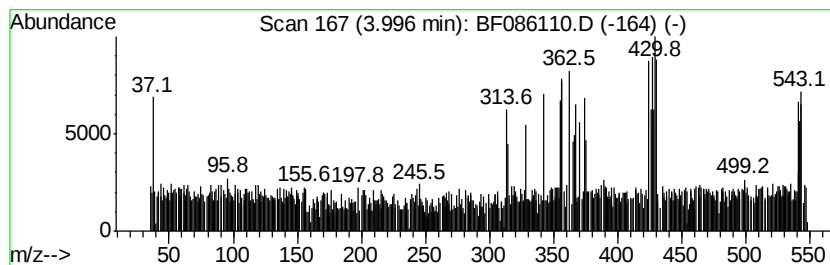
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown4.00 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.00	77.65 ng	2736270	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N,5-Bis[3,4-dichlorophenyl]-3,5-...	540	C27H20Cl4N4	333399-40-9	30
2		Nickel, bis[1,2-diphenyl-1,2-eth...	542	C28H20NiS4	028984-20-5	10
3		Ferrocene, 1,1'-bis(2,3,4,5-tetr...	542	C24H14F8FeO2	055494-18-3	10
4		3-Bromo-1',4-dihydroxy-2,3'-dime...	542	C27H27BrO7	101459-33-0	9
5		Anthra[2,1,9-def:6,5,10-d'e'f']d...	542	C36H18N2O4	000128-65-4	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086110.D
Acq On : 6 Apr 2016 21:01
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Sample : H2074-01
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB01A

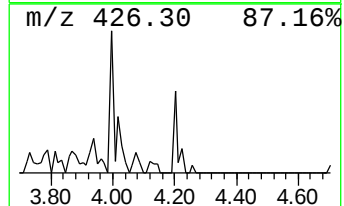
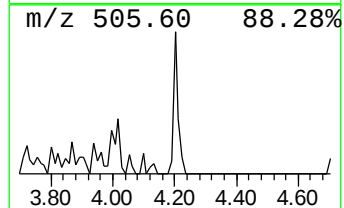
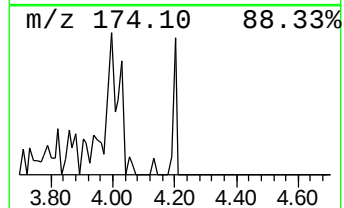
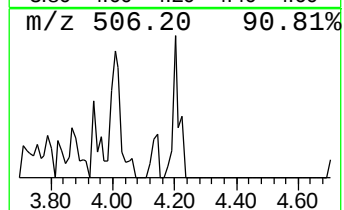
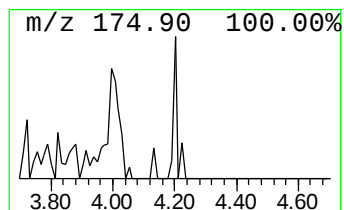
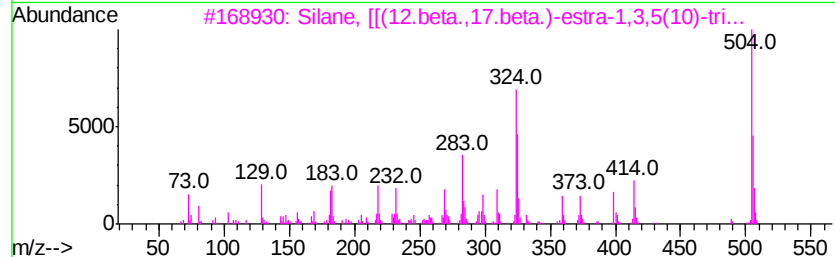
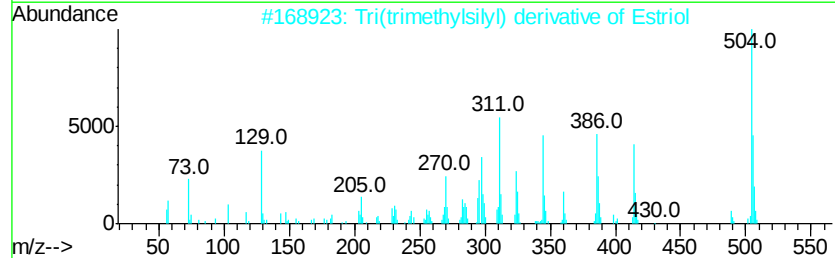
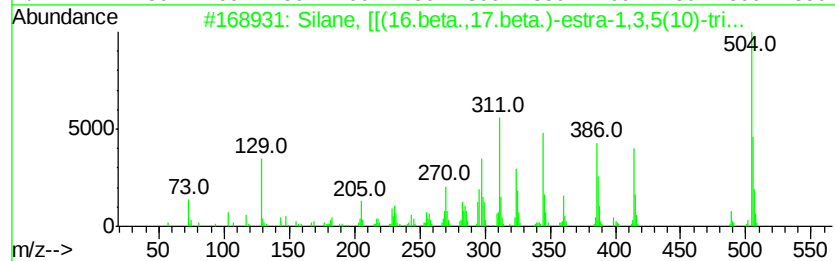
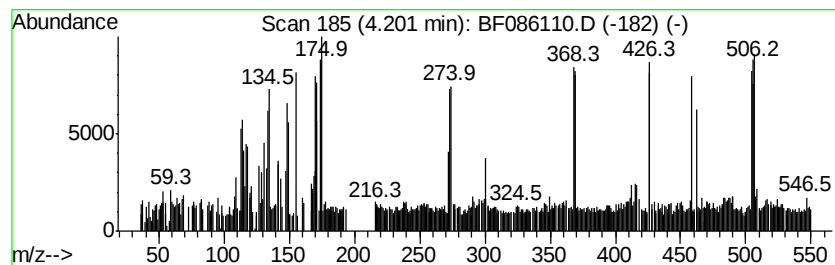
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown4.20 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	19.07 ng	671872	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, [[(16.beta.,17.beta.)-es...	504	C27H48O3Si3	057305-21-2	25
2		Tri(trimethylsilyl) derivative o...	504	C27H48O3Si3	018888-17-0	11
3		Silane, [[(12.beta.,17.beta.)-es...	504	C27H48O3Si3	069597-47-3	10
4		Silane, [[(17.beta.)-estra-1,3,5...	504	C27H48O3Si3	051497-42-8	10
5		Nickel, bis(1,3-diphenyl-1,3-pro...	504	C30H22NiO4	014405-47-1	9



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ALS Vial : 21 Sample Multiplier: 1

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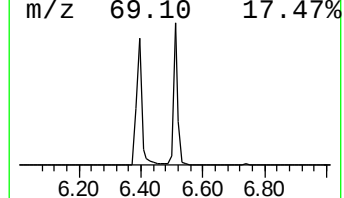
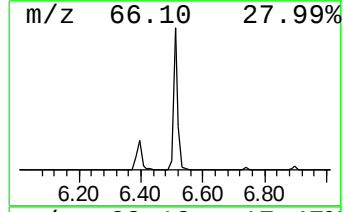
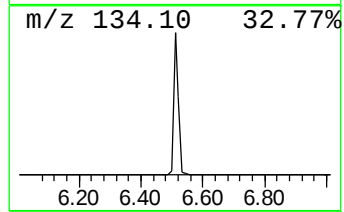
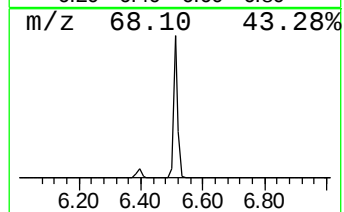
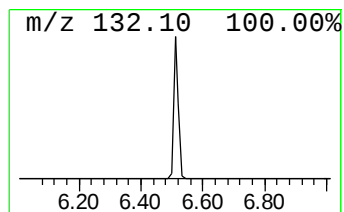
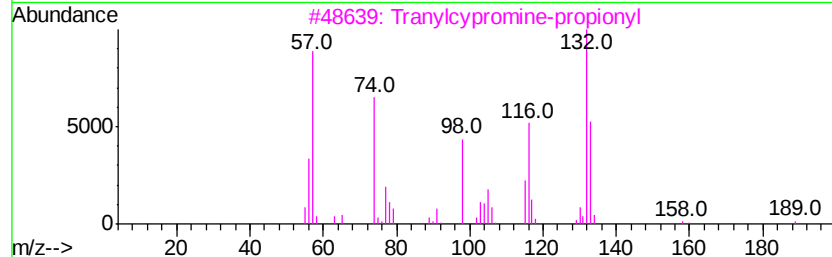
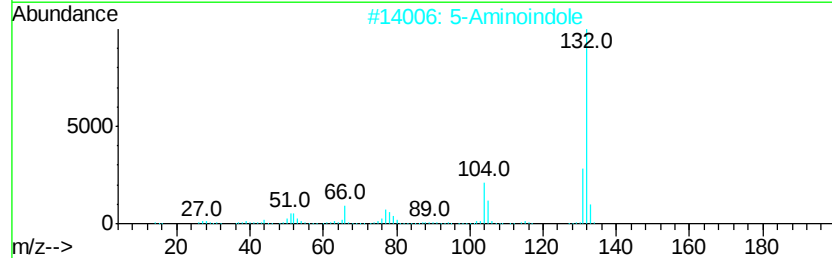
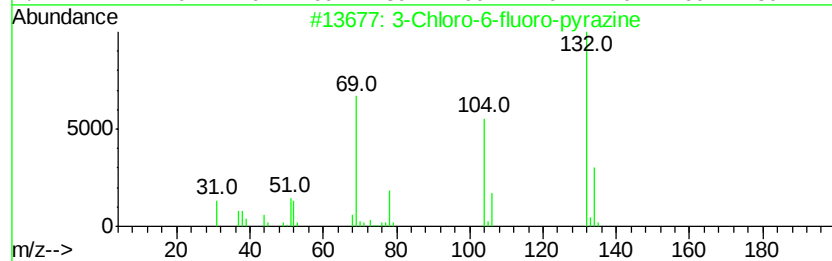
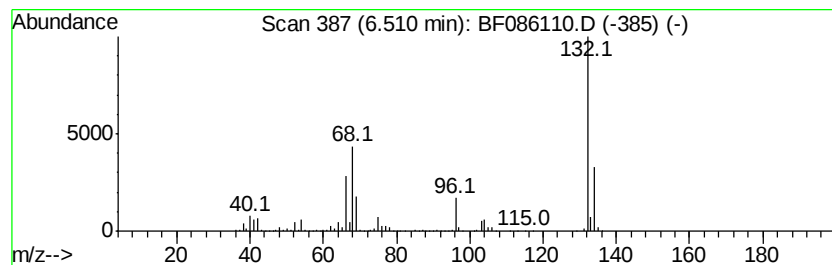
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown6.51 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	64.13 ng	2259950	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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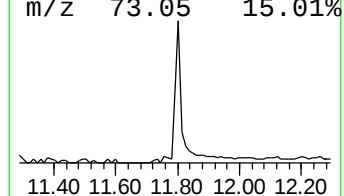
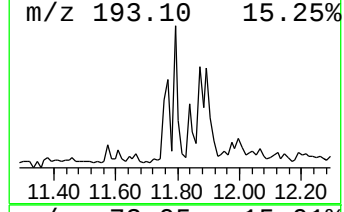
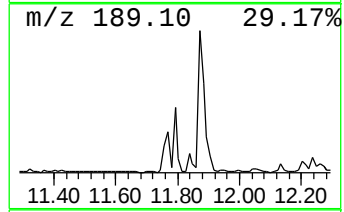
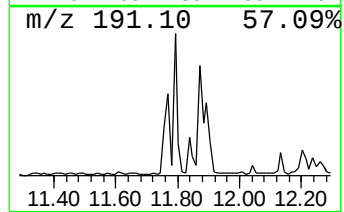
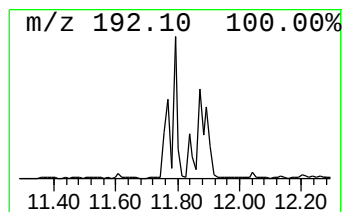
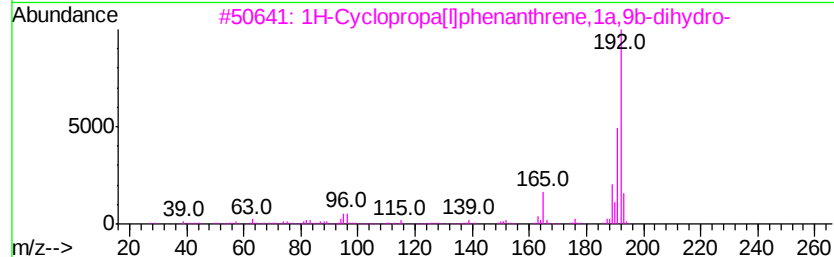
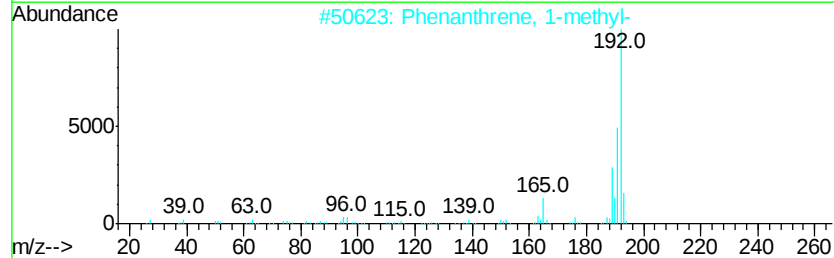
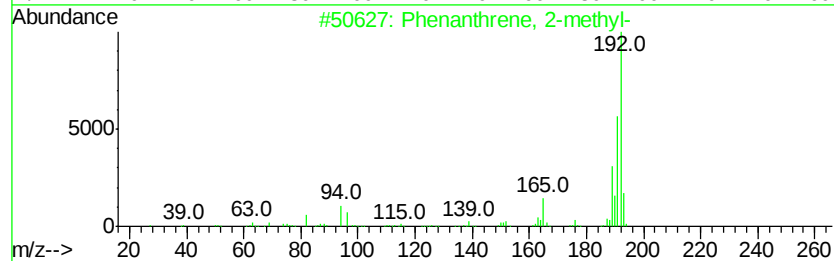
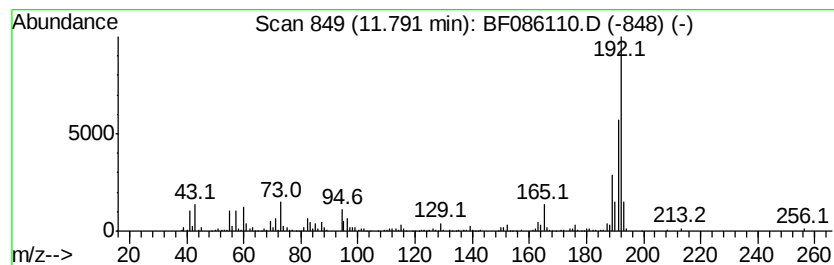
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Phenanthrene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	8.83 ng	373071	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



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Sample : H2074-01
Misc :
ALS Vial : 21 Sample Multiplier: 1

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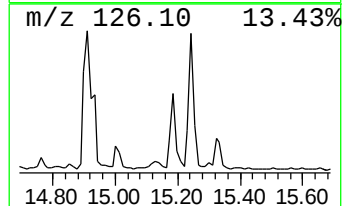
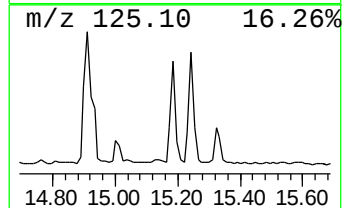
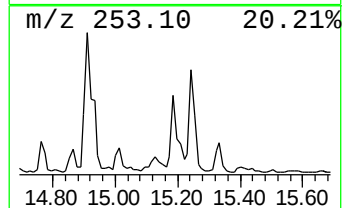
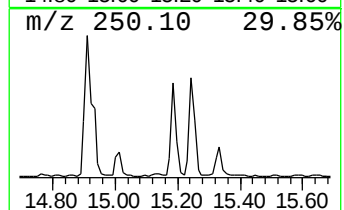
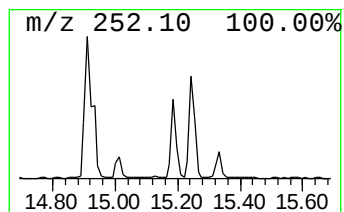
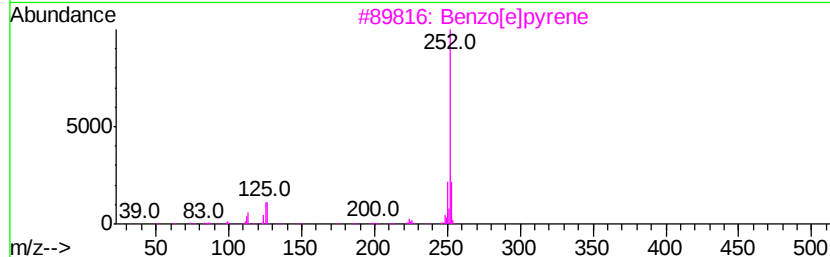
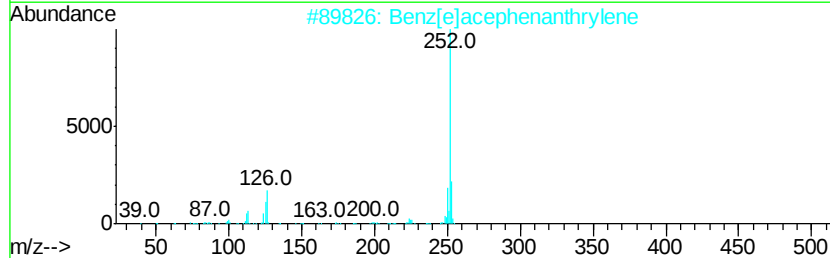
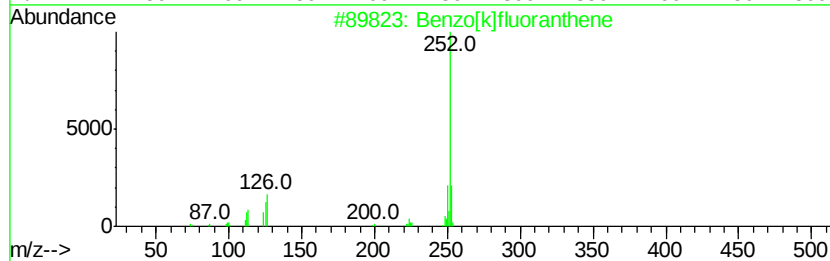
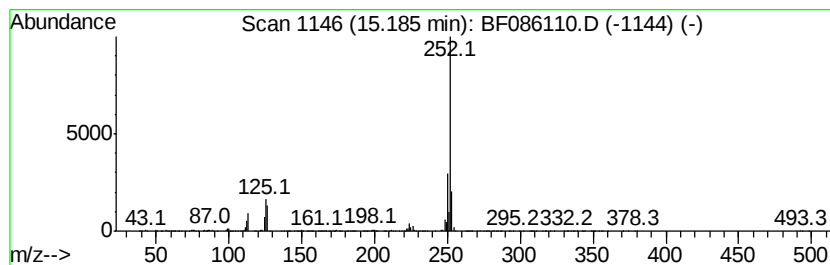
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzo[e]pyrene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	9.00 ng	337275	Perylene-d12	15.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	99
2			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
3			Benzo[e]pyrene	252	C20H12	000192-97-2	98
4			Perylene	252	C20H12	000198-55-0	95
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95



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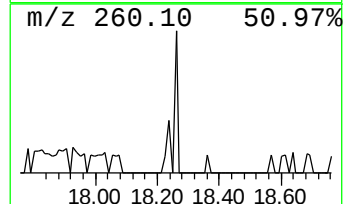
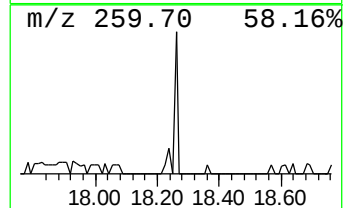
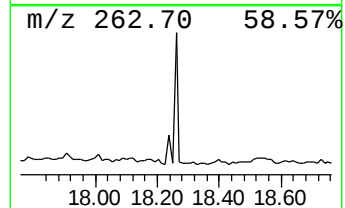
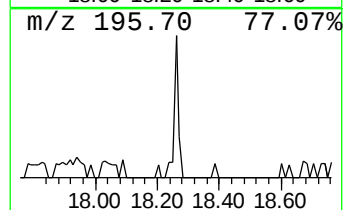
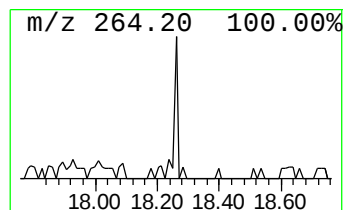
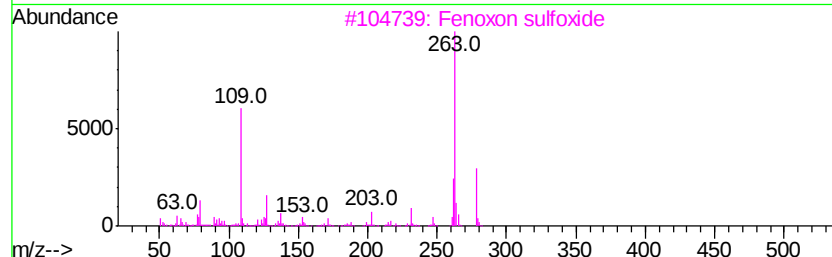
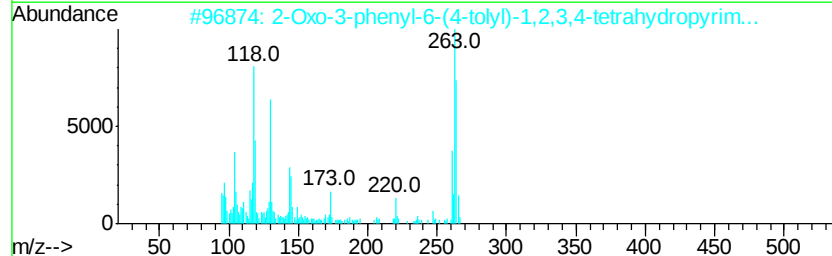
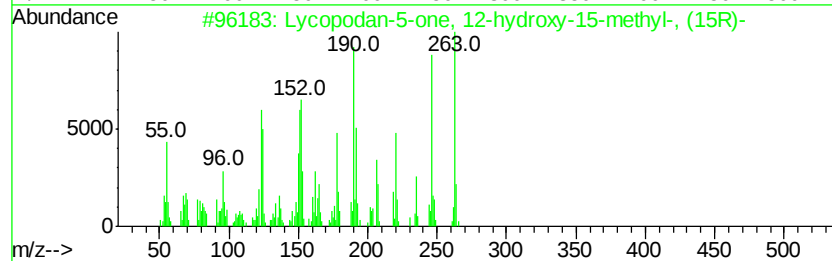
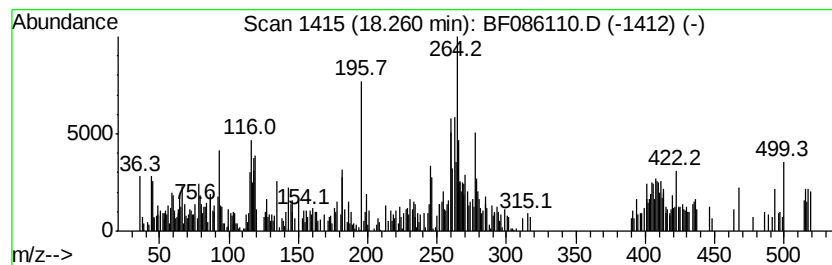
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Lycopodan-5-one, 12-hydroxy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.26	9.52 ng	356965	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Lycopodan-5-one, 12-hydroxy-15-methyl-, (15R)-	263	C16H25NO2	006900-92-1	56
2		2-Oxo-3-phenyl-6-(4-tolyl)-1,2,3,4-tetrahydropyrimidin-5(1H)-one	264	C17H16N2O	1000070-55-1	22
3		Fenoxon sulfoxide	278	C10H15O5PS	006552-13-2	11
4		Quinoline-4-carboxylic acid, 2-(4-tert-butylphenyl)-	263	C17H13NO2	020389-04-2	11
5		3,5-di-tert-Butyl-4-hydroxyphenylboronic acid	278	C17H26O3	020170-32-5	11



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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					#	RT	Resp	Conc
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unknown3.92	3.92	11.7	ng	412636	1	6.74	704793	20.0
unknown4.00	4.00	77.7	ng	2736270	1	6.74	704793	20.0
unknown4.20	4.20	19.1	ng	671872	1	6.74	704793	20.0
unknown6.51	6.51	64.1	ng	2259950	1	6.74	704793	20.0
Phenanthrene, 2-m...	11.79	8.8	ng	373071	4	11.27	845231	20.0
Benzo[e]pyrene	15.19	9.0	ng	337275	6	15.30	749647	20.0
Lycopodan-5-one, ...	18.26	9.5	ng	356965	6	15.30	749647	20.0