

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
 Data File : BF093227.D
 Acq On : 28 Feb 2017 1:42
 Operator : SJ/MA
 Sample : I1997-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-W7-BASE

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-BF013017.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	3.687	137	140	144	rVV	632983	1072225	14.21%	0.640%
2	4.339	194	197	205	rBV	487221	865180	11.47%	0.516%
3	4.979	249	253	257	rBV	1115290	1764680	23.39%	1.053%
4	5.253	274	277	279	rVV	889048	1223615	16.22%	0.730%
5	5.367	284	287	290	rBV	1547663	2707243	35.88%	1.615%
6	5.424	290	292	295	rVB	3587334	3743583	49.62%	2.233%
7	5.630	305	310	313	rVV	1279681	1707725	22.63%	1.019%
8	5.699	314	316	318	rVB	911526	822270	10.90%	0.491%
9	6.339	367	372	373	rBV2	609116	758187	10.05%	0.452%
10	6.407	376	378	381	rVB2	421549	663319	8.79%	0.396%
11	6.476	381	384	386	rBV	3534773	4250531	56.34%	2.536%
12	6.590	392	394	397	rVB	3288520	3760570	49.84%	2.244%
13	6.647	397	399	402	rBV	1084186	1563359	20.72%	0.933%
14	6.819	412	414	418	rBV2	1140738	1201474	15.92%	0.717%
15	6.899	418	421	422	rBV2	368838	537236	7.12%	0.321%
16	6.979	426	428	430	rBV	3169441	3021546	40.05%	1.803%
17	7.082	431	437	440	rBV4	313441	779139	10.33%	0.465%
18	7.390	462	464	466	rVV2	2476461	2595791	34.41%	1.549%
19	7.619	479	484	486	rBV4	266843	557979	7.40%	0.333%
20	7.848	500	504	507	rBV4	298616	733288	9.72%	0.437%
21	8.042	517	521	523	rBV4	163141	493384	6.54%	0.294%
22	8.133	523	529	531	rBV2	2379308	4163651	55.19%	2.484%
23	8.693	573	578	581	rVV3	459059	1013229	13.43%	0.605%
24	8.819	587	589	591	rVV	1541409	1249347	16.56%	0.745%
25	8.922	595	598	599	rVV	771066	965240	12.79%	0.576%
26	9.185	617	621	623	rVB	3683764	3775186	50.04%	2.252%
27	9.276	626	629	632	rBV2	403688	843699	11.18%	0.503%
28	9.448	641	644	645	rBV	358990	513509	6.81%	0.306%
29	9.516	648	650	651	rVV	758006	714335	9.47%	0.426%
30	9.539	651	652	654	rVV2	254789	478991	6.35%	0.286%
31	9.585	654	656	658	rVV	1585826	1713890	22.72%	1.023%
32	9.722	665	668	670	rBV	1691516	1767809	23.43%	1.055%
33	9.791	672	674	676	rVB	644800	593160	7.86%	0.354%
34	9.859	676	680	681	rBV2	1292926	1518363	20.12%	0.906%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.893	681	683	685	rVV	1228303	1164678	15.44%	0.695%
36	10.019	691	694	696	rBV3	336723	549801	7.29%	0.328%
37	10.065	696	698	699	rVV	568125	632122	8.38%	0.377%
38	10.282	713	717	719	rBV4	354349	831455	11.02%	0.496%
39	10.408	726	728	731	rVV	1231153	1304024	17.28%	0.778%
40	10.511	736	737	740	rVV2	442543	622718	8.25%	0.372%
41	10.648	747	749	751	rVV2	1506261	1828952	24.24%	1.091%
42	10.762	757	759	763	rVB2	838385	1358492	18.01%	0.810%
43	10.956	773	776	777	rBV	422062	593310	7.86%	0.354%
44	11.105	785	789	791	rBV2	432805	642514	8.52%	0.383%
45	11.208	795	798	799	rBV2	582324	662287	8.78%	0.395%
46	11.242	799	801	804	rVB	643499	728953	9.66%	0.435%
47	11.379	808	813	815	rBV	4768083	6589082	87.33%	3.931%
48	11.425	815	817	818	rVV	2178542	1739125	23.05%	1.038%
49	11.585	829	831	833	rBV	499039	565194	7.49%	0.337%
50	11.642	834	836	837	rVB	787846	721106	9.56%	0.430%
51	11.848	852	854	856	rVV	1288788	1817325	24.09%	1.084%
52	11.882	856	857	859	rVV	1579746	1424957	18.89%	0.850%
53	11.928	859	861	862	rVV	824607	881676	11.69%	0.526%
54	11.962	862	864	868	rVV	1934050	3394992	45.00%	2.026%
55	12.054	869	872	874	rVV3	402632	926501	12.28%	0.553%
56	12.088	874	875	877	rVV2	549508	645507	8.56%	0.385%
57	12.145	878	880	882	rVB2	788087	1018994	13.51%	0.608%
58	12.294	891	893	895	rVB3	519338	616078	8.17%	0.368%
59	12.408	902	903	904	rBV	714714	597859	7.92%	0.357%
60	12.442	904	906	909	rVB3	783196	1346171	17.84%	0.803%
61	12.499	909	911	913	rVB2	483598	800207	10.61%	0.477%
62	12.579	915	918	919	rBV	4126874	4933851	65.39%	2.944%
63	12.602	919	920	922	rVB2	1187266	1228446	16.28%	0.733%
64	12.808	935	938	940	rBV	4295757	7544709	100.00%	4.501%
65	12.854	940	942	944	rVB2	442682	623778	8.27%	0.372%
66	12.934	947	949	953	rVB2	1784115	2646813	35.08%	1.579%
67	13.014	953	956	960	rBV3	884221	1726101	22.88%	1.030%
68	13.128	962	966	968	rVB	1573775	2299528	30.48%	1.372%
69	13.197	970	972	973	rBV2	967560	1105539	14.65%	0.660%
70	13.231	973	975	979	rBV	1276469	1480966	19.63%	0.884%
71	13.322	981	983	985	rVV	1337956	1631373	21.62%	0.973%

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 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
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72	13.357	985	986	987	rVB	1387102	954475	12.65%	0.569%
73	13.642	1006	1011	1014	rBV2	479399	1305739	17.31%	0.779%
74	13.745	1017	1020	1021	rBV2	507316	898693	11.91%	0.536%
75	13.768	1021	1022	1024	rVV	580500	579367	7.68%	0.346%
76	13.997	1039	1042	1044	rBV2	2971718	6141993	81.41%	3.664%
77	14.031	1044	1045	1049	rVV	2951690	3034736	40.22%	1.811%
78	14.111	1049	1052	1054	rVV2	941761	1307640	17.33%	0.780%
79	14.157	1054	1056	1060	rVB3	534344	917110	12.16%	0.547%
80	14.385	1071	1076	1078	rBV	1184749	2251076	29.84%	1.343%
81	14.465	1081	1083	1086	rBV3	655524	1343241	17.80%	0.801%
82	14.580	1091	1093	1096	rVB2	664204	776512	10.29%	0.463%
83	14.877	1117	1119	1121	rBV	350709	492875	6.53%	0.294%
84	15.037	1129	1133	1138	rBV	2391678	6244158	82.76%	3.725%
85	15.128	1140	1141	1144	rVB	784078	941513	12.48%	0.562%
86	15.334	1155	1159	1161	rBV	2553768	4322941	57.30%	2.579%
87	15.403	1161	1165	1167	rVV	2850047	5448352	72.21%	3.251%
88	15.471	1170	1171	1174	rVB2	894133	1364732	18.09%	0.814%
89	15.745	1193	1195	1197	rBV	386751	597296	7.92%	0.356%
90	15.780	1197	1198	1201	rVB	453453	516354	6.84%	0.308%
91	15.985	1214	1216	1220	rVB2	564881	946436	12.54%	0.565%
92	16.866	1288	1293	1296	rBV	1428297	3382922	44.84%	2.018%
93	17.014	1303	1306	1308	rBV	290327	557514	7.39%	0.333%
94	17.311	1326	1332	1335	rBV	1534438	4315497	57.20%	2.575%
95	17.517	1346	1350	1355	rVB	495139	1250820	16.58%	0.746%
96	18.089	1397	1400	1407	rVB7	248850	623429	8.26%	0.372%
97	18.397	1424	1427	1437	rVB5	204735	964930	12.79%	0.576%
98	19.174	1492	1495	1498	rVB4	206560	511441	6.78%	0.305%
99	19.289	1498	1505	1507	rBV4	239583	918314	12.17%	0.548%
100	19.449	1515	1519	1525	rVB	387426	1343957	17.81%	0.802%

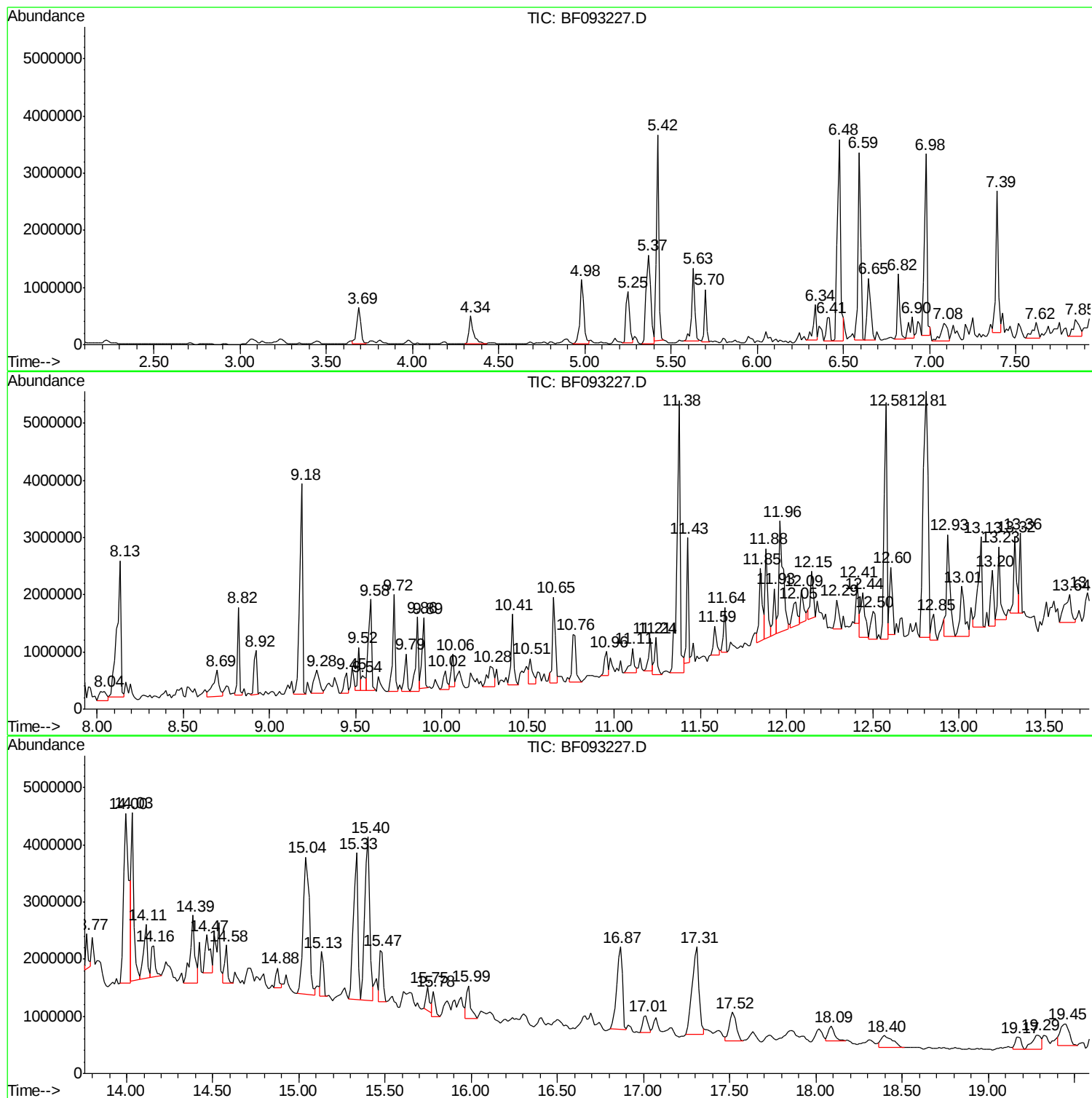
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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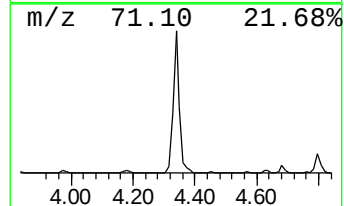
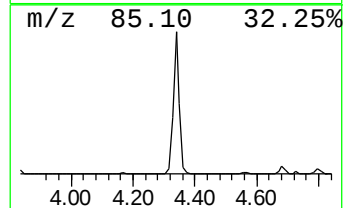
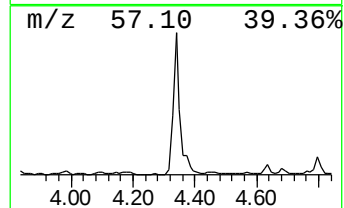
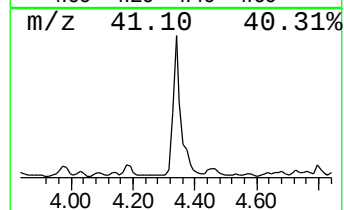
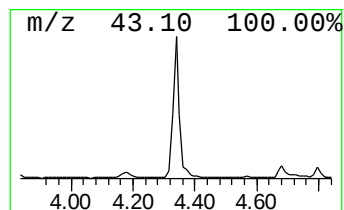
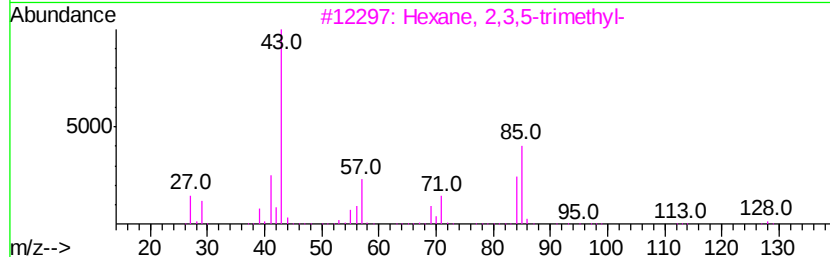
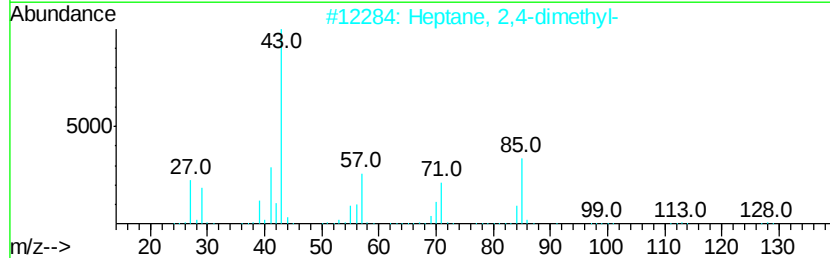
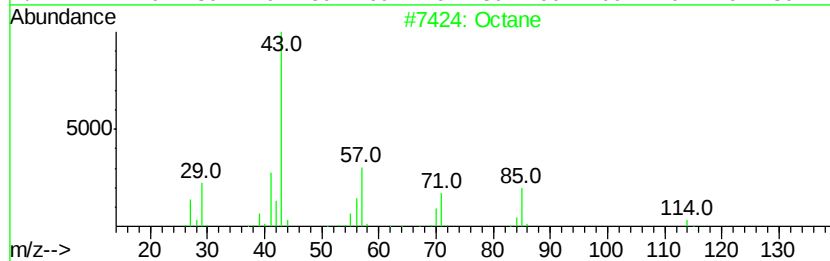
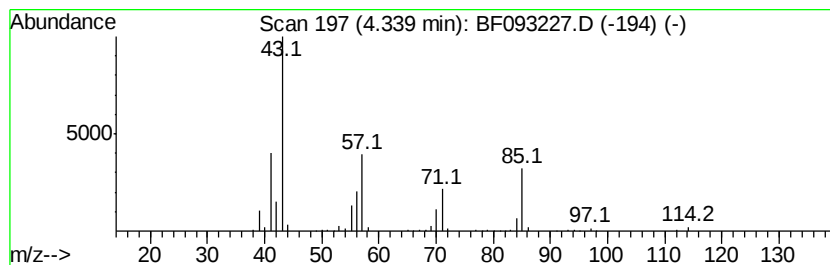
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Octane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.34	14.40 ng	865180	1,4-Dichlorobenzene-d4	6.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane		114	C8H18	000111-65-9	91
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	80
3	Hexane, 2,3,5-trimethyl-		128	C9H20	001069-53-0	64
4	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	59
5	Hexane, 3-ethyl-		114	C8H18	000619-99-8	52



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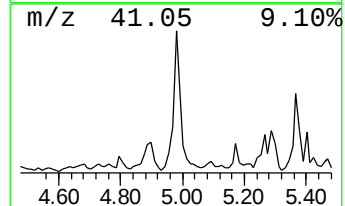
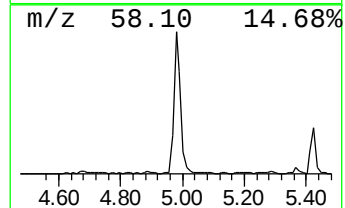
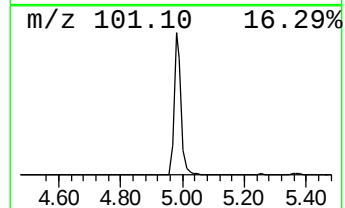
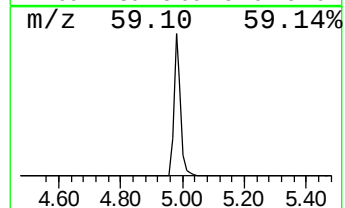
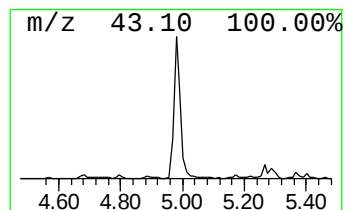
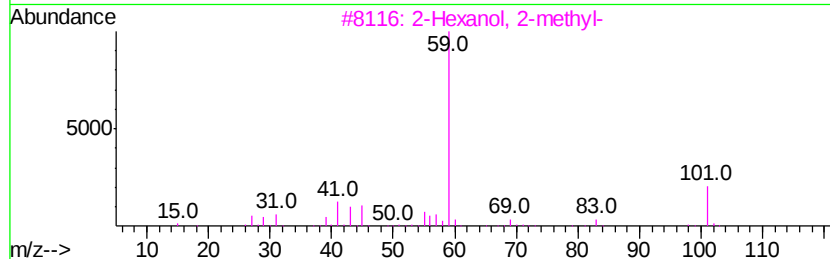
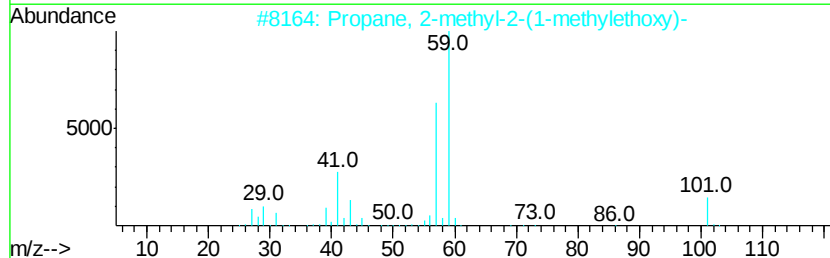
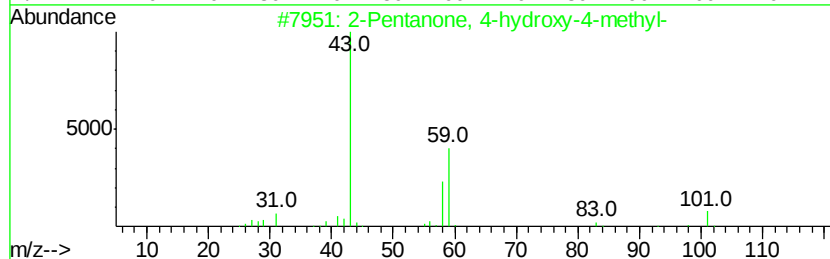
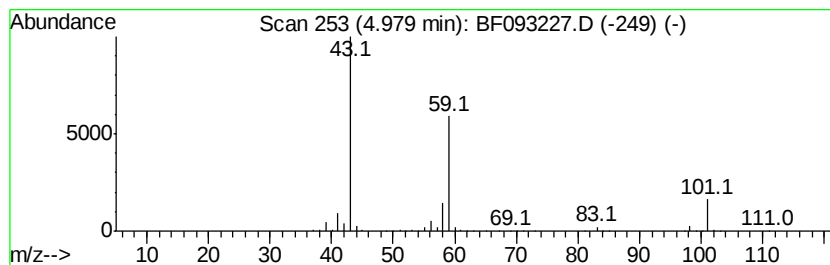
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	29.38 ng	1764680	1,4-Dichlorobenzene-d4	6.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O		017348-59-3	33
3	2-Hexanol, 2-methyl-	116	C7H16O		000625-23-0	33
4	3-Hexanol, 4-methyl-	116	C7H16O		000615-29-2	33
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2		001116-98-9	23



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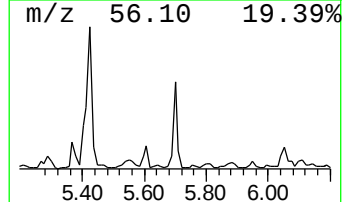
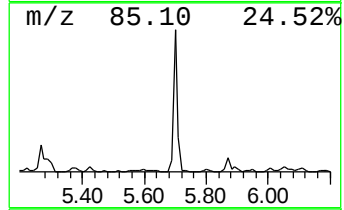
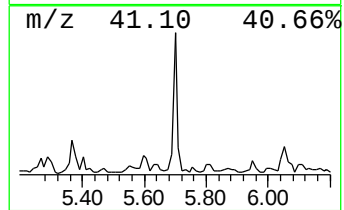
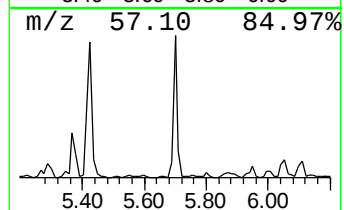
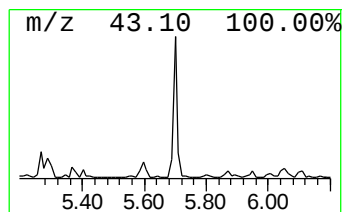
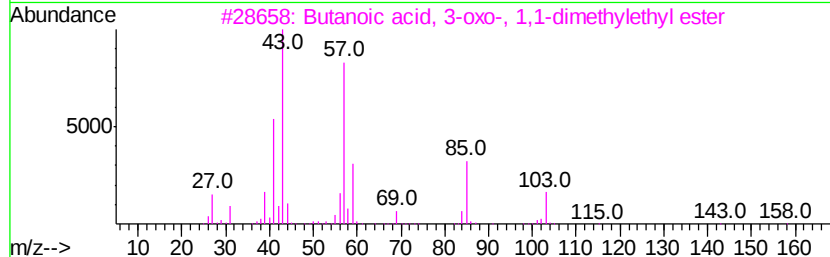
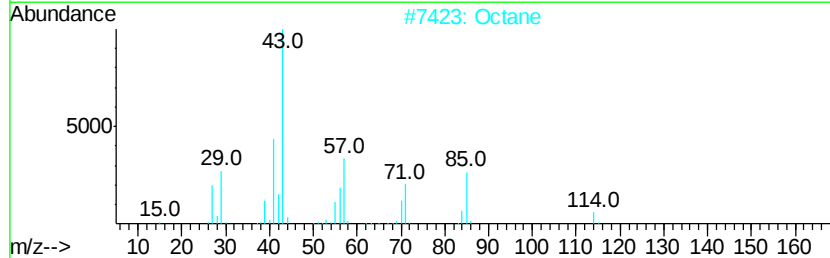
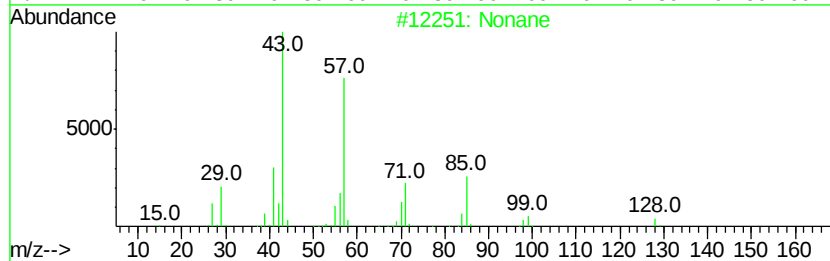
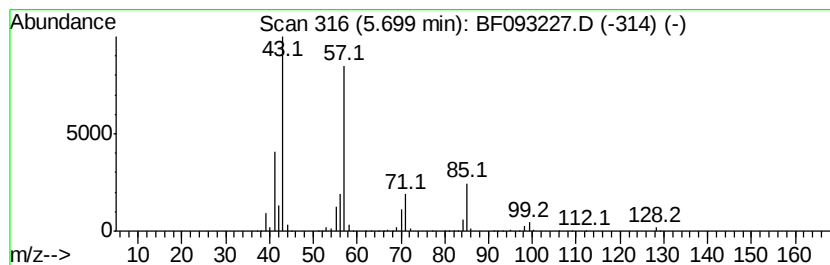
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Nonane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	13.69 ng	822270	1,4-Dichlorobenzene-d4	6.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	90
2	Octane		114	C8H18	000111-65-9	64
3	Butanoic acid, 3-oxo-, 1,1-dimet...		158	C8H14O3	001694-31-1	59
4	Decane, 2,6,7-trimethyl-		184	C13H28	062108-25-2	59
5	Heptane, 2,3-dimethyl-		128	C9H20	003074-71-3	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
Operator : SJ/MA
Sample : I1997-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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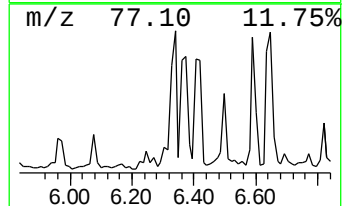
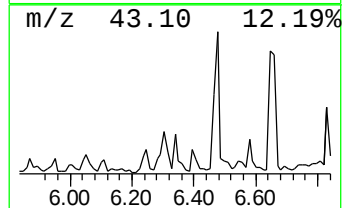
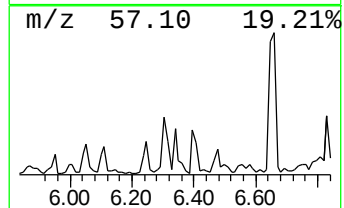
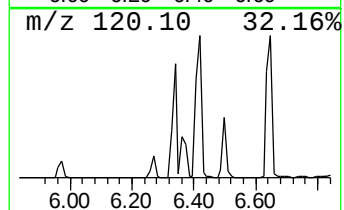
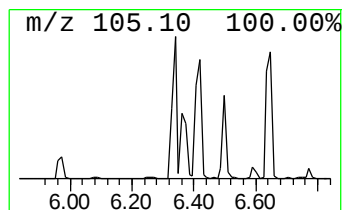
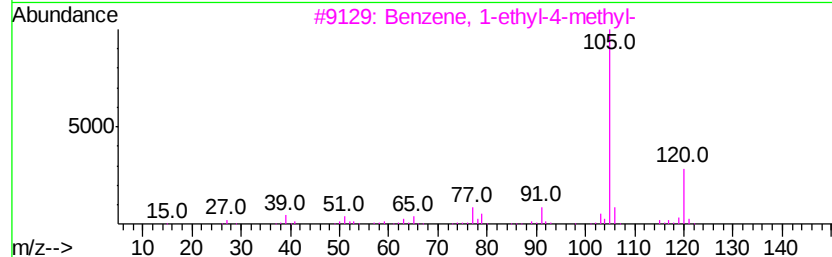
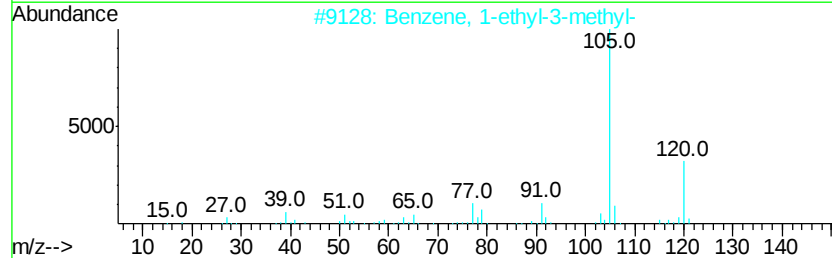
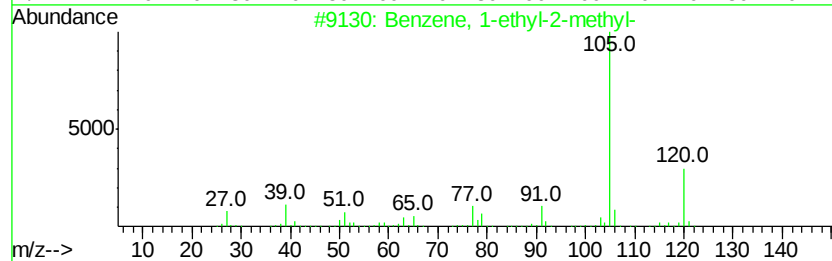
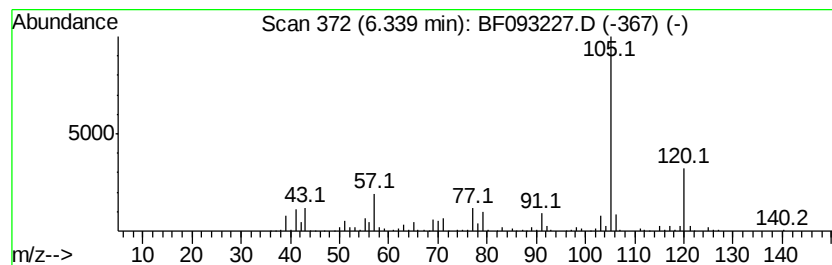
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	12.62 ng	758187	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
Operator : SJ/MA
Sample : I1997-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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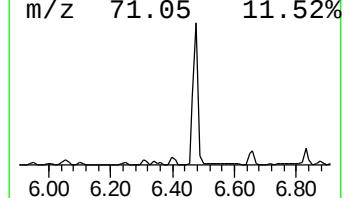
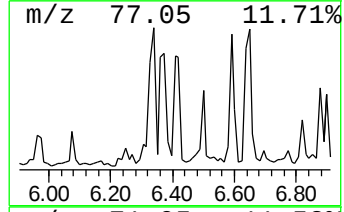
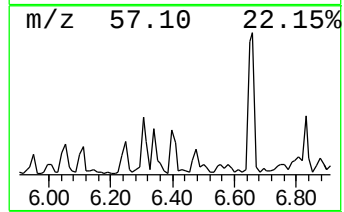
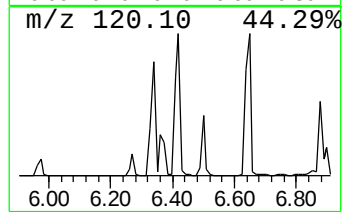
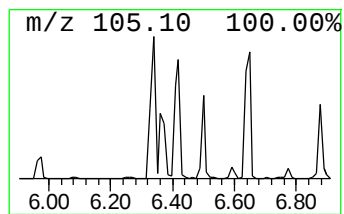
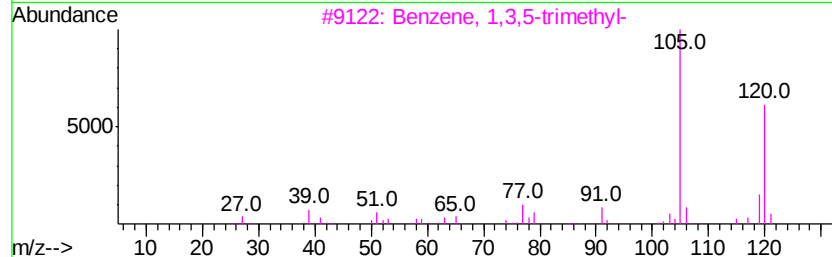
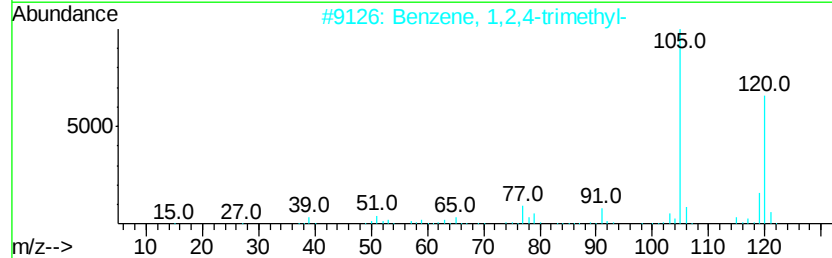
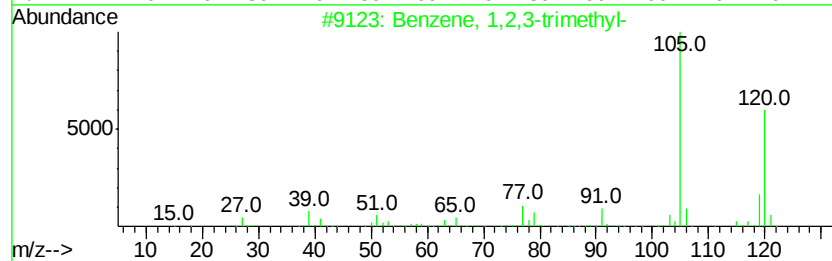
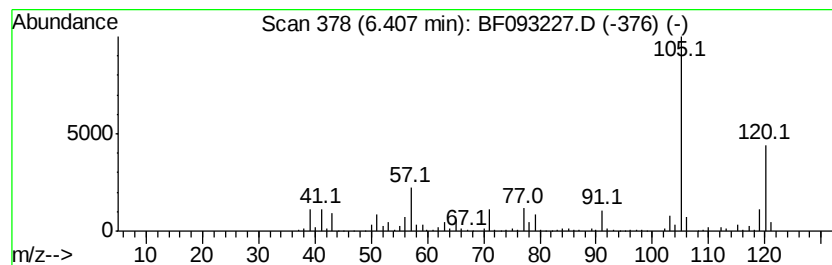
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	11.04 ng	663319	1,4-Dichlorobenzene-d4	6.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
Operator : SJ/MA
Sample : I1997-04
Misc :
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Instrument :
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ClientSampleId :
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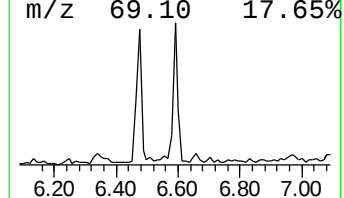
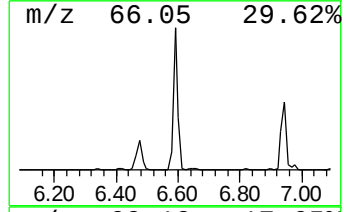
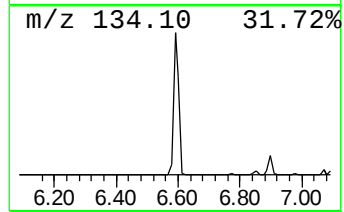
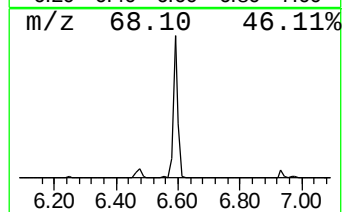
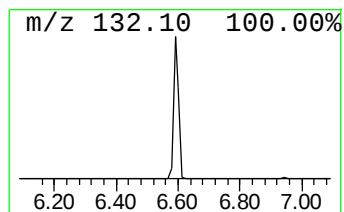
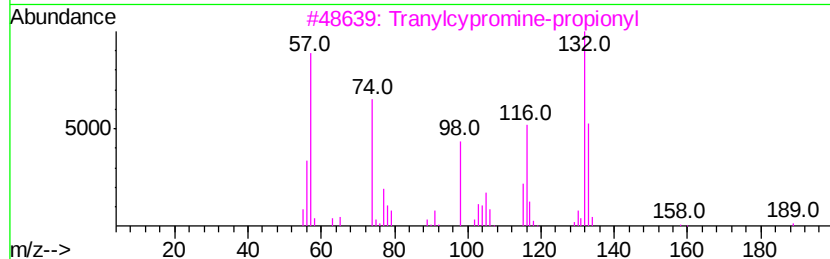
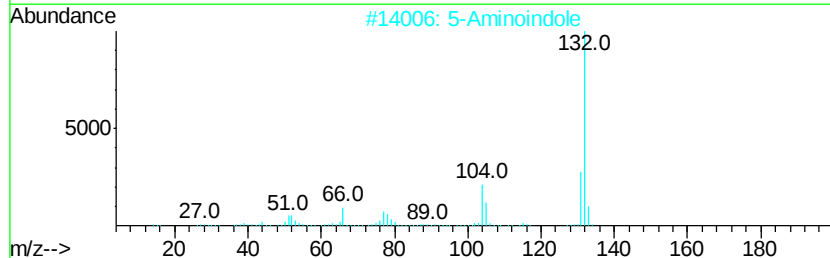
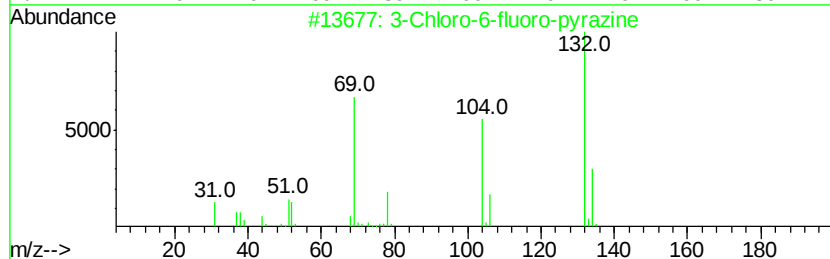
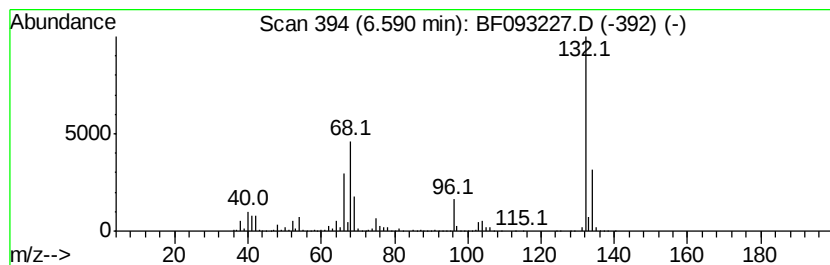
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown6.59 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	62.60 ng	3760570	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
Operator : SJ/MA
Sample : I1997-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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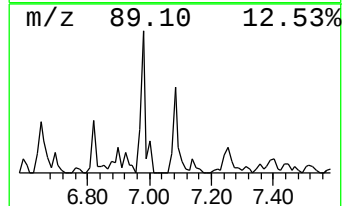
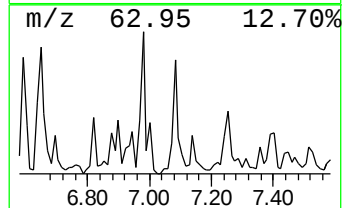
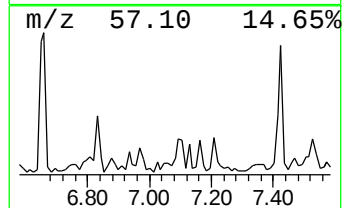
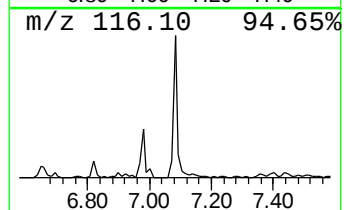
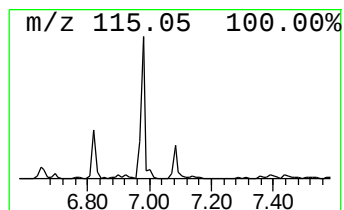
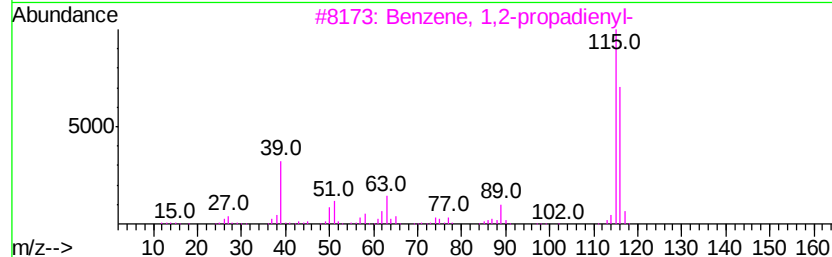
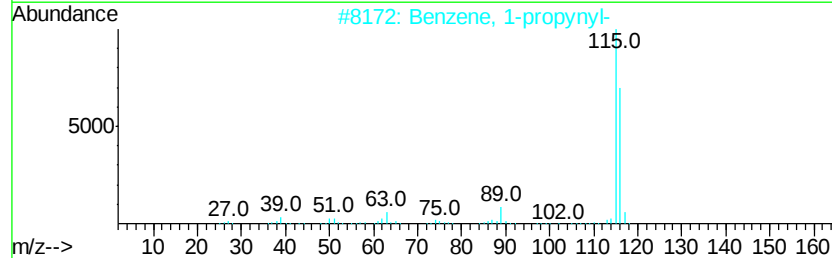
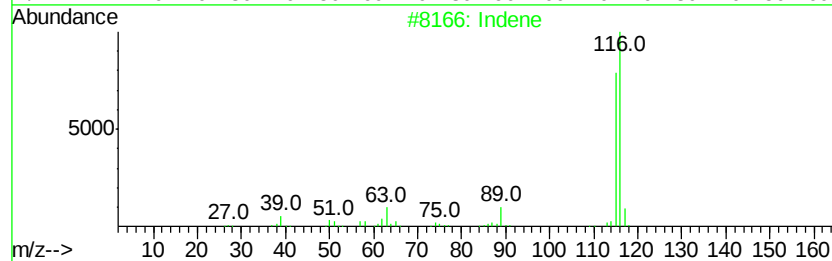
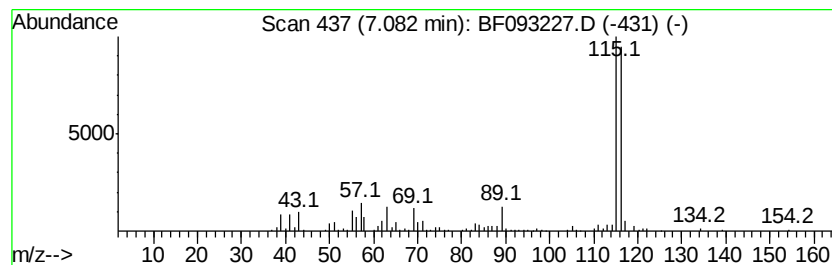
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Indene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	12.97 ng	779139	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	91
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	87
3		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	80
4		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	80
5		1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
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Sample : I1997-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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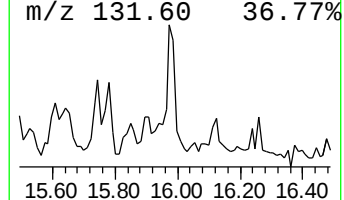
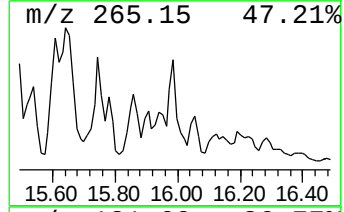
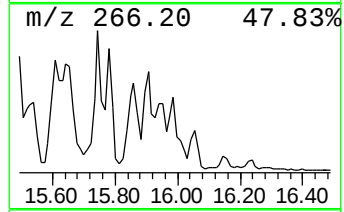
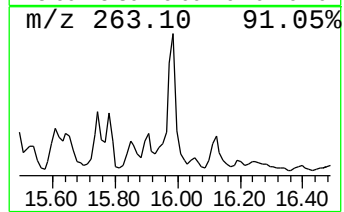
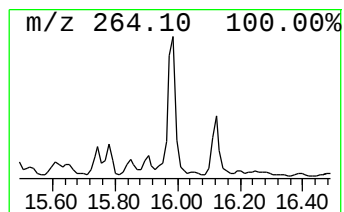
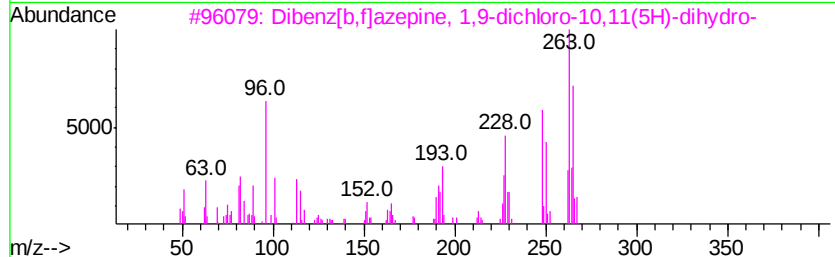
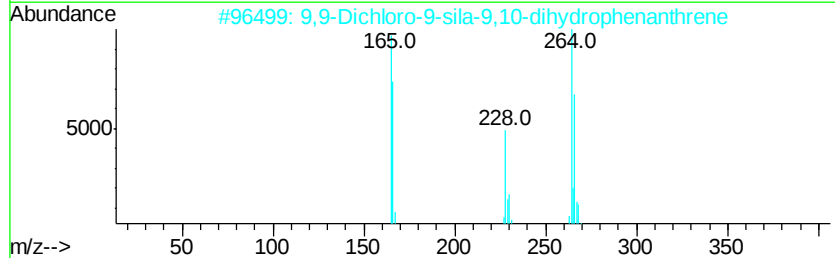
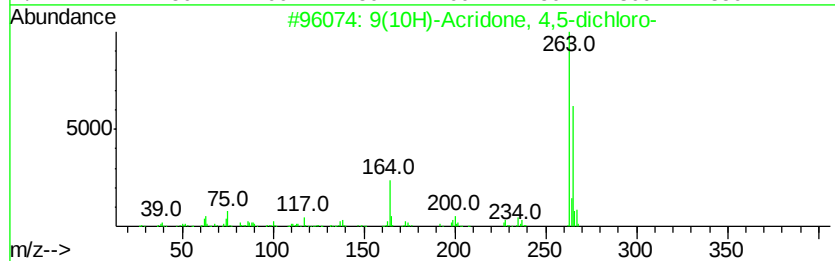
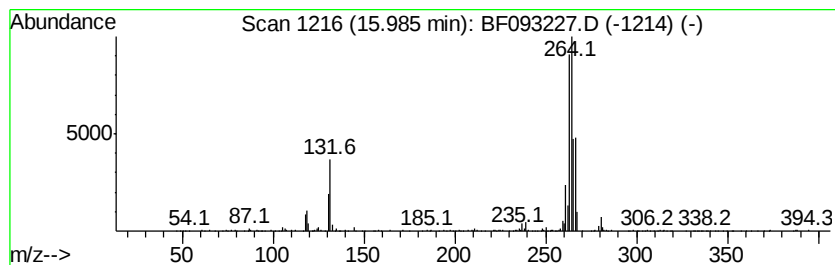
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown15.99 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	13.87 ng	946436	Perylene-d12	15.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9(10H)-Acridone, 4,5-dichloro-			263	C13H7Cl2NO	1000163-26-5	35
2	9,9-Dichloro-9-sila-9,10-dihydro...			264	C13H10Cl2Si	076470-26-3	32
3	Dibenz[b,f]azepine, 1,9-dichloro...			263	C14H11Cl2N	107517-53-3	32
4	Quinoline, 2-[2-(4-chlorophenyl)]...			265	C17H12ClN	005392-19-8	32
5	Boroxin, ethyldiphenyl-			264	C14H15B3O3	1000151-82-0	28



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
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Misc :
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ClientSampleId :
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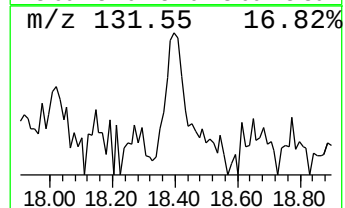
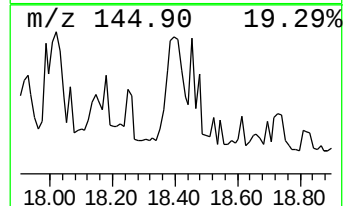
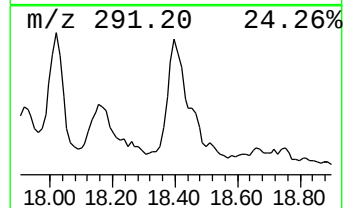
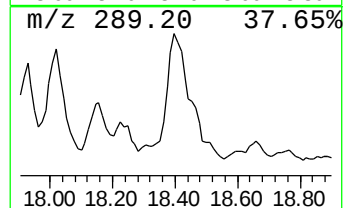
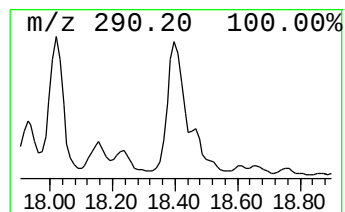
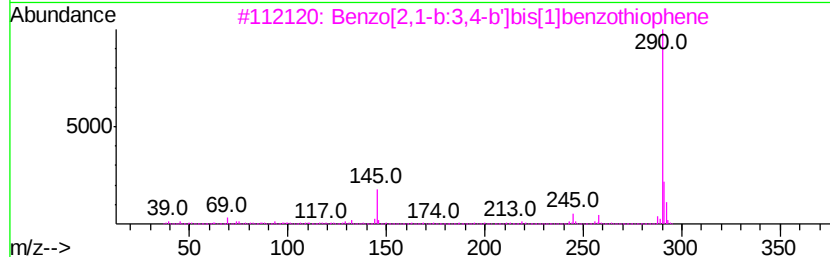
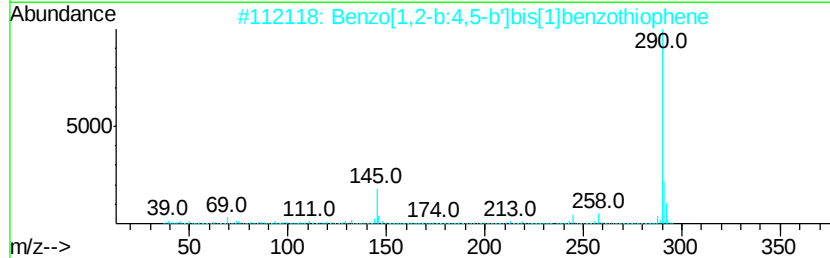
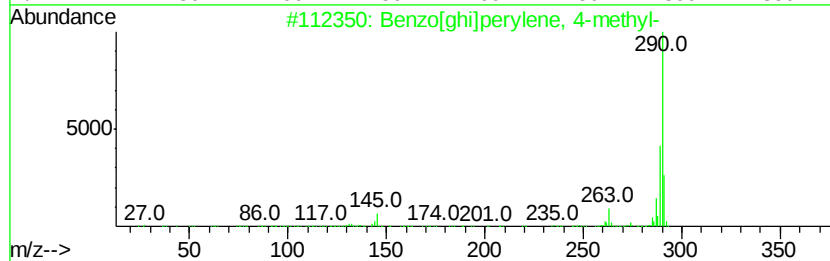
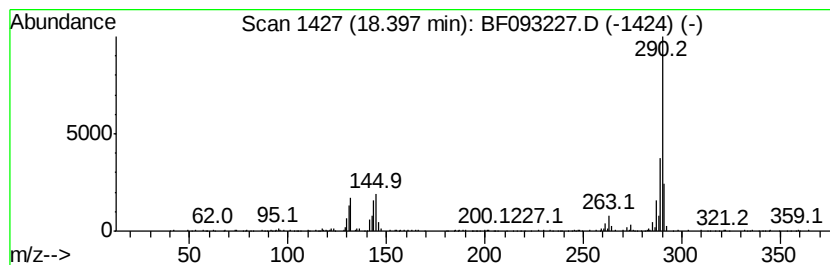
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[ghi]perylene, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	14.14 ng	964930	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[ghi]perylene, 4-methyl-	290	C23H14	019224-38-5	89
2		Benzo[1,2-b:4,5-b']bis[1]benzoth...	290	C18H10S2	000241-34-9	52
3		Benzo[2,1-b:3,4-b']bis[1]benzoth...	290	C18H10S2	000222-24-2	50
4		Benzo[1,2-b:3,4-b']bis[1]benzoth...	290	C18H10S2	000198-89-0	50
5		11H-Indeno[1,2-b]-1,2,5-oxadiaz...	290	C15H6N4OS	302602-04-6	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
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ALS Vial : 16 Sample Multiplier: 1

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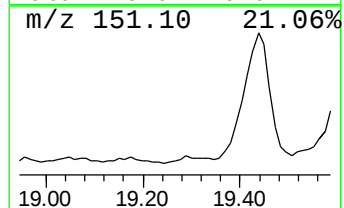
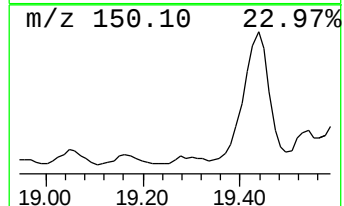
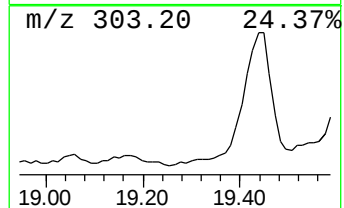
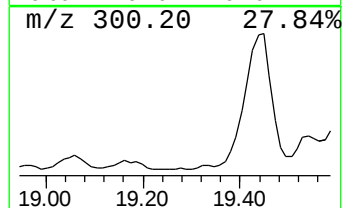
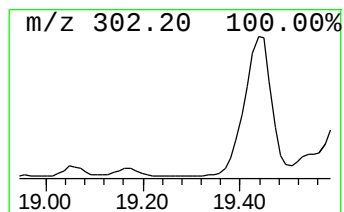
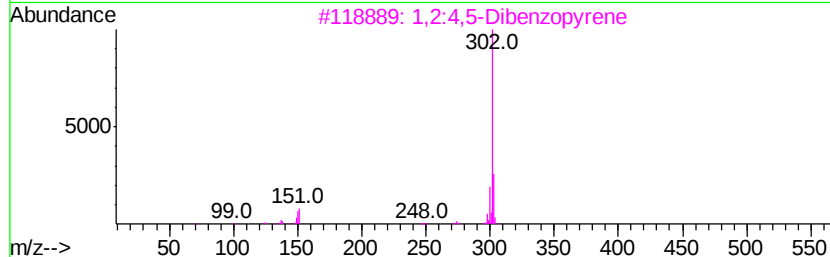
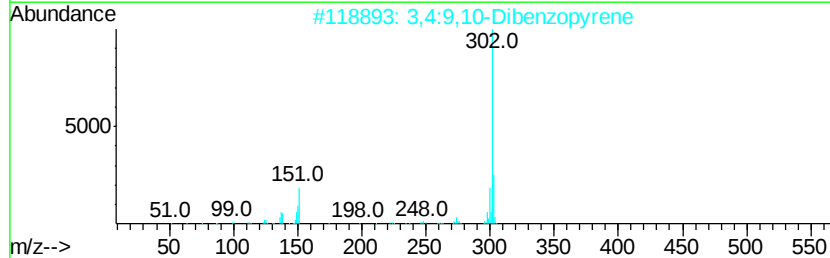
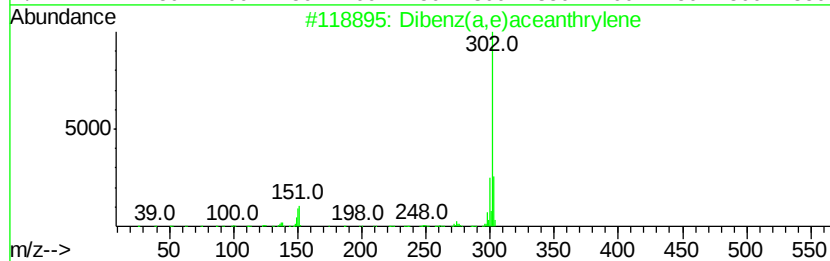
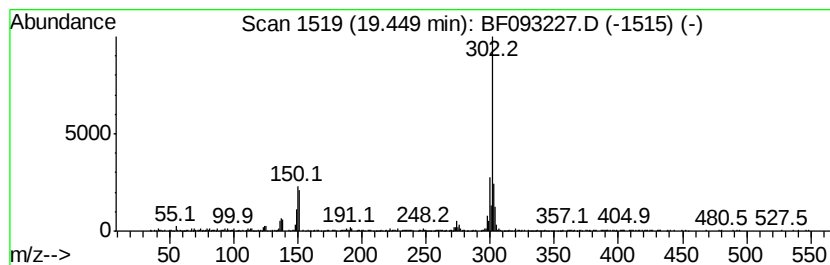
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Dibenz(a,e)aceanthrylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.45	19.70 ng	1343960	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	98
2		3,4:9,10-Dibenzopyrene	302	C24H14	000189-55-9	97
3		1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	96
4		3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	93
5		Indeno[1,2,3-fg]naphthacene	302	C24H14	000203-11-2	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093227.D
Acq On : 28 Feb 2017 1:42
Operator : SJ/MA
Sample : I1997-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-W7-BASE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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
Octane	4.34	14.4	ng	865180	1	6.82	1201470	20.0
2-Pentanone, 4-hy...	4.98	29.4	ng	1764680	1	6.82	1201470	20.0
Nonane	5.70	13.7	ng	822270	1	6.82	1201470	20.0
Benzene, 1-ethyl-...	6.34	12.6	ng	758187	1	6.82	1201470	20.0
Benzene, 1,2,3-tr...	6.41	11.0	ng	663319	1	6.82	1201470	20.0
unknown6.59	6.59	62.6	ng	3760570	1	6.82	1201470	20.0
Indene	7.08	13.0	ng	779139	1	6.82	1201470	20.0
unknown15.99	15.99	13.9	ng	946436	6	15.45	1364730	20.0
Benzo[ghi]perylene...	18.40	14.1	ng	964930	6	15.45	1364730	20.0
Friedelan-3-one	19.29	13.5	ng	918314	6	15.45	1364730	20.0
Dibenz(a,e)aceant...	19.45	19.7	ng	1343960	6	15.45	1364730	20.0