

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
 Data File : BG031590.D
 Acq On : 15 Dec 2017 11:30
 Operator : SJ/JU
 Sample : I6830-04
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MAP-5-E(12-13)

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_G\METHODS\8270-BG121117.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	5.658	396	403	416	rBV	1192027	2097284	17.31%	0.943%
2	6.351	513	521	528	rBV5	133264	388467	3.21%	0.175%
3	6.463	534	540	548	rVB3	81741	184697	1.52%	0.083%
4	6.633	565	569	573	rVB	174661	239864	1.98%	0.108%
5	6.739	582	587	598	rVB4	128365	323698	2.67%	0.146%
6	6.898	609	614	620	rVB4	109640	199538	1.65%	0.090%
7	7.268	665	677	690	rVV	1199005	2758170	22.77%	1.240%
8	7.409	696	701	707	rVB4	147119	289404	2.39%	0.130%
9	7.467	707	711	715	rBV3	148629	234636	1.94%	0.106%
10	7.632	733	739	746	rVB	1255414	2289929	18.90%	1.030%
11	7.808	765	769	775	rVB4	152692	272310	2.25%	0.122%
12	7.914	779	787	792	rBV3	210110	469604	3.88%	0.211%
13	7.996	792	801	807	rBV7	418971	1229294	10.15%	0.553%
14	8.061	807	812	816	rBV	797913	1157322	9.55%	0.520%
15	8.161	823	829	834	rVB4	329838	672544	5.55%	0.302%
16	8.231	834	841	843	rBV6	124347	279743	2.31%	0.126%
17	8.272	843	848	853	rVB3	481499	763715	6.30%	0.343%
18	8.343	853	860	863	rBV2	293184	631866	5.22%	0.284%
19	8.425	868	874	883	rVB	1087940	2296991	18.96%	1.033%
20	8.537	885	893	897	rVB6	126826	326101	2.69%	0.147%
21	8.590	897	902	903	rBV2	219960	278222	2.30%	0.125%
22	8.654	910	913	922	rVB3	408596	923118	7.62%	0.415%
23	8.737	922	927	930	rBV2	203836	376495	3.11%	0.169%
24	8.872	946	950	956	rBV7	133448	269616	2.23%	0.121%
25	8.936	956	961	966	rBV2	700864	1455124	12.01%	0.654%
26	9.083	982	986	995	rVB8	214081	584813	4.83%	0.263%
27	9.201	995	1006	1012	rBV2	1464086	3415626	28.20%	1.536%
28	9.271	1012	1018	1029	rVB2	768107	2020169	16.68%	0.908%
29	9.359	1030	1033	1037	rVB3	356765	537562	4.44%	0.242%
30	9.412	1037	1042	1043	rBV2	549371	710132	5.86%	0.319%
31	9.447	1044	1048	1057	rVB6	709643	1700586	14.04%	0.765%
32	9.524	1057	1061	1062	rBV2	242624	295708	2.44%	0.133%
33	9.565	1063	1068	1075	rVB5	753330	1824240	15.06%	0.820%
34	9.677	1076	1087	1092	rBV9	594862	1874902	15.48%	0.843%

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35	9.735	1092	1097	1102	rBV2	1795797	3072243	25.36%	1.382%
36	9.829	1108	1113	1123	rVB3	1176154	2759708	22.78%	1.241%
37	9.988	1130	1140	1144	rBV5	944905	2344453	19.35%	1.054%
38	10.088	1151	1157	1164	rBV5	1711392	3794109	31.32%	1.706%
39	10.147	1164	1167	1170	rVV2	317690	455260	3.76%	0.205%
40	10.305	1190	1194	1197	rBV4	246449	495920	4.09%	0.223%
41	10.370	1198	1205	1215	rVB6	847564	2286504	18.88%	1.028%
42	10.487	1219	1225	1234	rBV6	1238026	3074692	25.38%	1.383%
43	10.605	1240	1245	1254	rBV2	1334592	3145441	25.97%	1.415%
44	10.687	1255	1259	1263	rBV4	490995	865773	7.15%	0.389%
45	10.769	1269	1273	1274	rBV2	367584	493130	4.07%	0.222%
46	10.875	1286	1291	1295	rBV6	615205	1494145	12.33%	0.672%
47	11.069	1318	1324	1329	rVB	3699988	6698216	55.30%	3.012%
48	11.128	1329	1334	1341	rBV3	1544648	2865498	23.66%	1.289%
49	11.292	1356	1362	1366	rBV5	665961	1545082	12.76%	0.695%
50	11.339	1366	1370	1375	rVB4	711248	1238167	10.22%	0.557%
51	11.404	1376	1381	1386	rVB4	1305629	2417326	19.96%	1.087%
52	11.463	1387	1391	1393	rBV4	490142	819592	6.77%	0.369%
53	11.574	1406	1410	1411	rBV4	596309	811781	6.70%	0.365%
54	11.686	1424	1429	1433	rBV3	717815	1196972	9.88%	0.538%
55	11.745	1435	1439	1445	rVV7	676325	1431693	11.82%	0.644%
56	11.933	1467	1471	1472	rBV2	772402	1028234	8.49%	0.462%
57	11.962	1472	1476	1480	rVV3	1743126	2757850	22.77%	1.240%
58	12.115	1496	1502	1507	rBV5	1191211	2713392	22.40%	1.220%
59	12.238	1520	1523	1528	rVB2	985981	1540349	12.72%	0.693%
60	12.362	1538	1544	1553	rBV10	1031222	3202040	26.43%	1.440%
61	12.650	1588	1593	1600	rBV4	1054580	2992121	24.70%	1.346%
62	13.267	1694	1698	1703	rVB4	1670940	2816108	23.25%	1.266%
63	13.355	1709	1713	1720	rVB3	2436053	4374339	36.11%	1.967%
64	13.531	1739	1743	1748	rBV5	1474043	3115988	25.72%	1.401%
65	13.619	1755	1758	1760	rBV4	835871	1150355	9.50%	0.517%
66	13.795	1784	1788	1793	rBV	2020950	3174822	26.21%	1.428%
67	14.166	1847	1851	1855	rBV	3599589	5130600	42.36%	2.307%
68	14.207	1855	1858	1863	rVB2	3132596	4218207	34.82%	1.897%
69	14.471	1898	1903	1911	rBV4	3748760	7606805	62.80%	3.421%
70	14.571	1916	1920	1925	rVB	4282003	6527030	53.88%	2.935%
71	14.818	1958	1962	1966	rBV5	1847105	2940352	24.27%	1.322%

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72	14.959	1982	1986	1991	rBV3	2062984	3014746	24.89%	1.356%
73	15.241	2031	2034	2037	rBV4	893359	1550549	12.80%	0.697%
74	15.358	2048	2054	2065	rBV	5549239	12113187	100.00%	5.447%
75	15.511	2075	2080	2085	rBV4	2783912	6498763	53.65%	2.923%
76	15.558	2085	2088	2095	rVB	3016910	4438127	36.64%	1.996%
77	15.781	2122	2126	2132	rVB3	3375580	5928870	48.95%	2.666%
78	15.916	2146	2149	2152	rBV	2247045	2708548	22.36%	1.218%
79	15.993	2158	2162	2167	rVB3	2425208	3446296	28.45%	1.550%
80	16.228	2199	2202	2210	rBV4	2767293	5180819	42.77%	2.330%
81	16.451	2237	2240	2252	rVB2	3425074	6361986	52.52%	2.861%
82	16.592	2259	2264	2268	rBV2	2510450	3922044	32.38%	1.764%
83	16.851	2303	2308	2313	rBV3	3493833	5577694	46.05%	2.508%
84	16.997	2329	2333	2335	rBV3	1295355	2072537	17.11%	0.932%
85	17.139	2354	2357	2361	rVB4	1202063	1551236	12.81%	0.698%
86	17.879	2480	2483	2489	rBV5	894205	1837862	15.17%	0.826%
87	18.202	2534	2538	2543	rBV2	896934	1471053	12.14%	0.662%
88	18.842	2644	2647	2650	rBV	500205	669323	5.53%	0.301%
89	19.089	2686	2689	2693	rVB	1306197	1656760	13.68%	0.745%
90	19.177	2701	2704	2709	rVB2	1108221	1439748	11.89%	0.647%
91	19.260	2713	2718	2722	rBV	2167774	3473229	28.67%	1.562%
92	19.389	2737	2740	2745	rBV	754915	1281192	10.58%	0.576%
93	19.442	2746	2749	2757	rVB	715309	1152273	9.51%	0.518%
94	19.747	2798	2801	2804	rBV2	391860	573999	4.74%	0.258%
95	19.782	2805	2807	2814	rVB	557480	833818	6.88%	0.375%
96	19.888	2821	2825	2830	rVB2	1019062	1614154	13.33%	0.726%
97	19.947	2831	2835	2840	rVB	1053015	1547303	12.77%	0.696%
98	20.064	2850	2855	2859	rVB	4227589	5445968	44.96%	2.449%
99	21.739	3134	3140	3155	rVB2	906470	1624122	13.41%	0.730%
100	25.012	3684	3697	3714	rBV2	474253	1417278	11.70%	0.637%

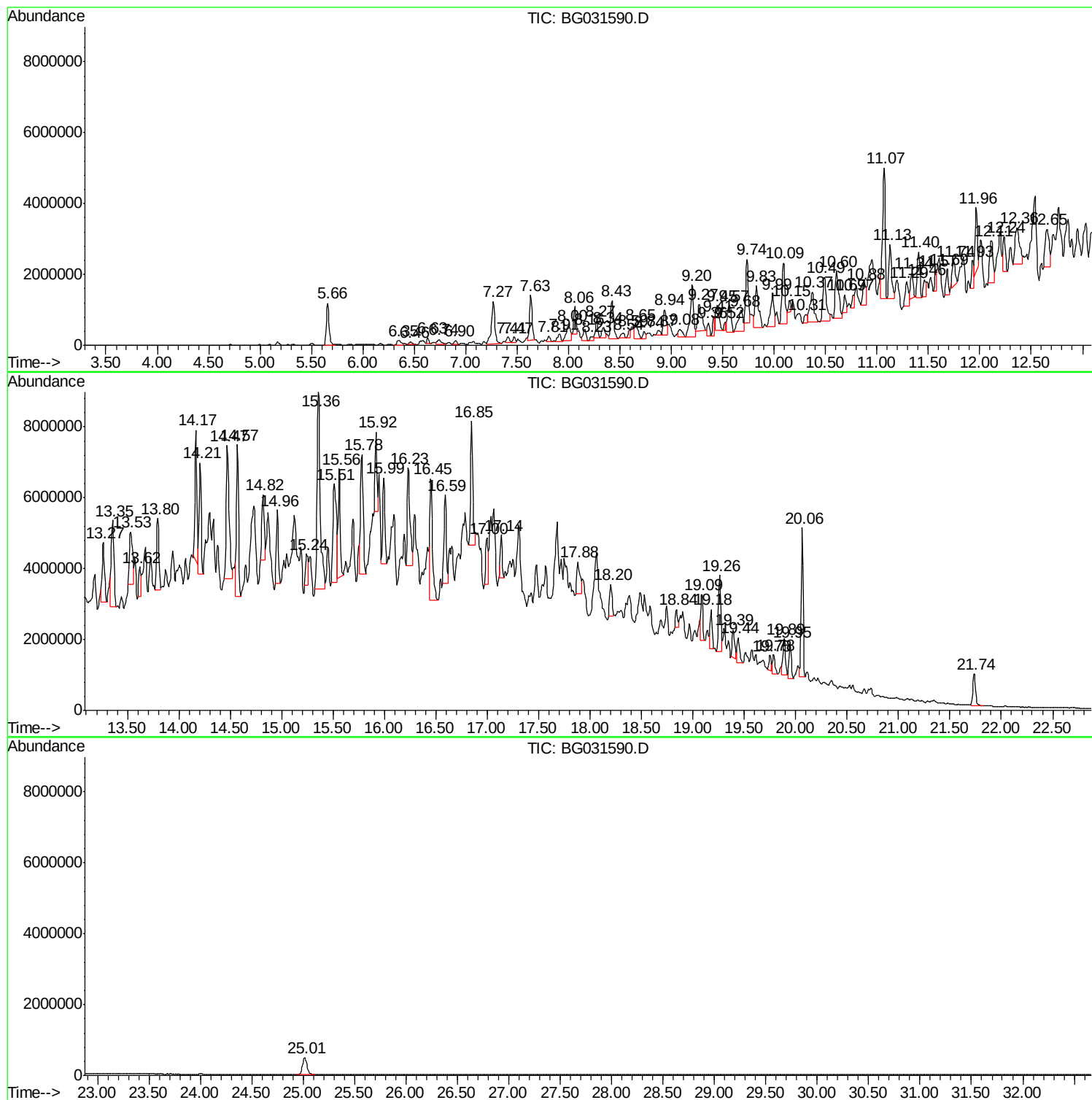
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Instrument :
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MAP-5-E(12-13)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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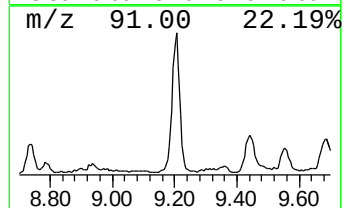
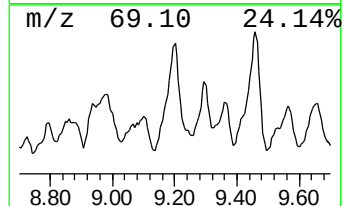
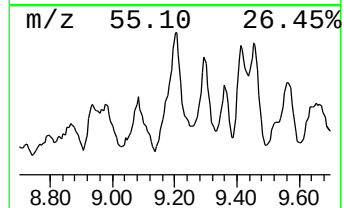
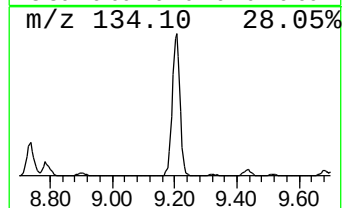
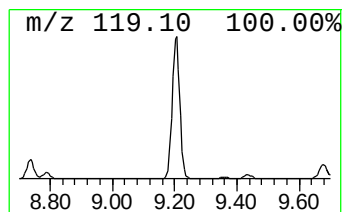
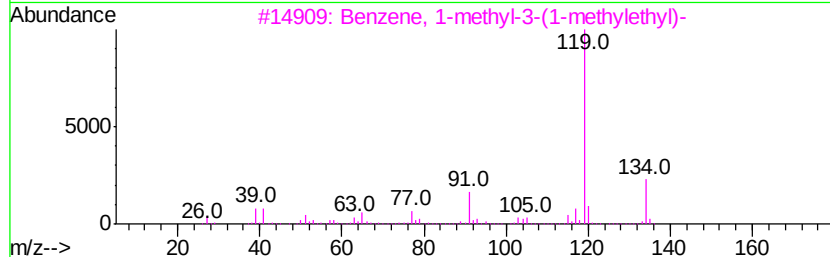
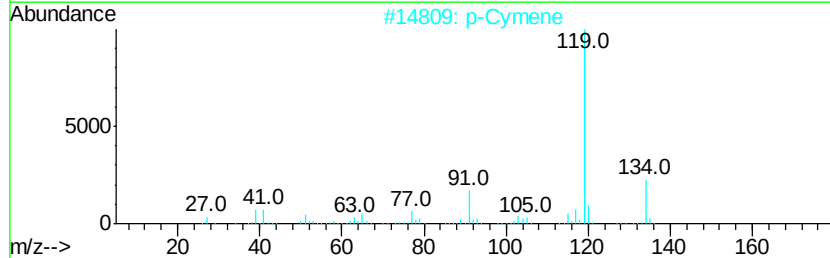
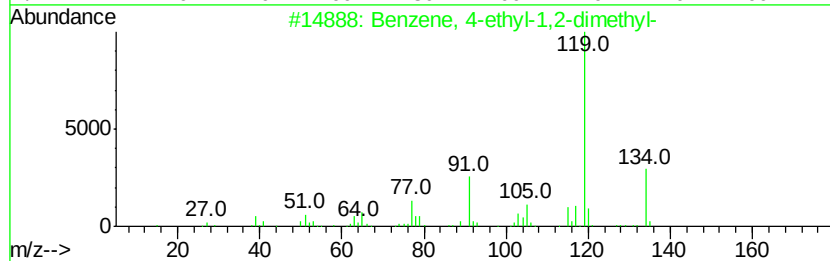
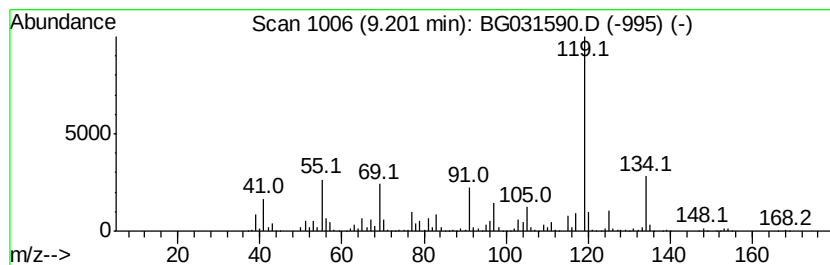
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.20	59.03 ng	3415630	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93



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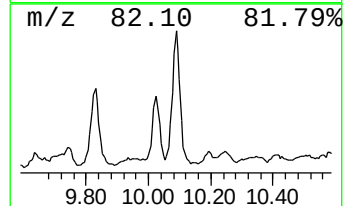
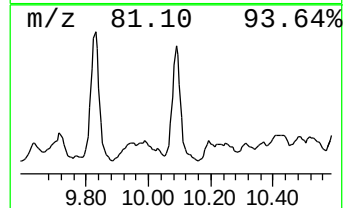
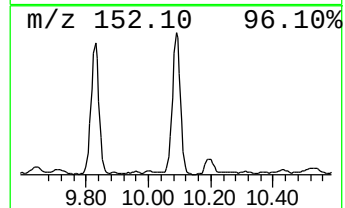
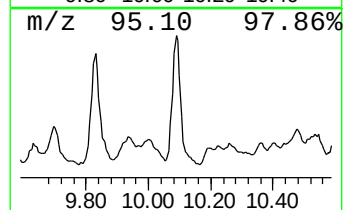
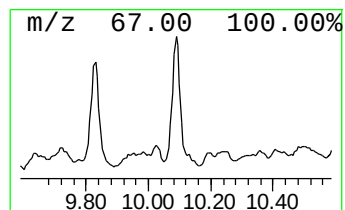
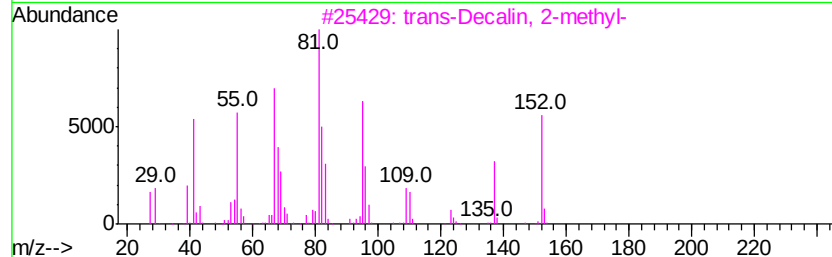
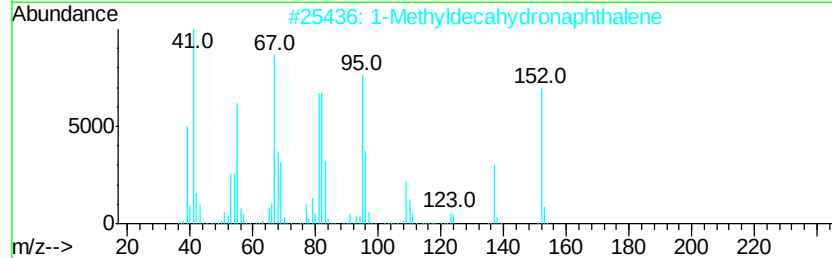
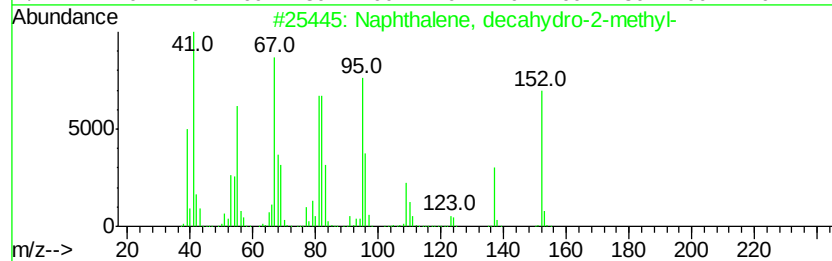
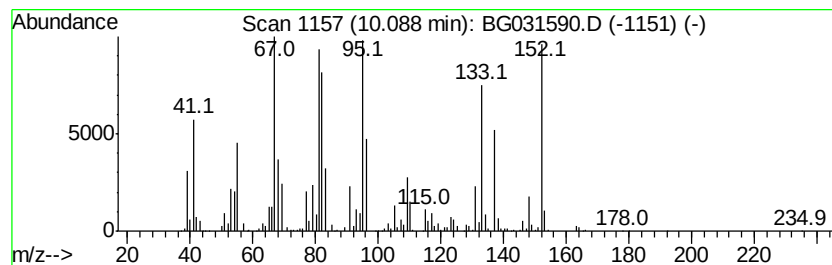
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Naphthalene, decahydro-2-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	50.79 ng	3794110	Naphthalene-d8	10.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	97
2		1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	97
3		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	62
4		cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	58
5		trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	55



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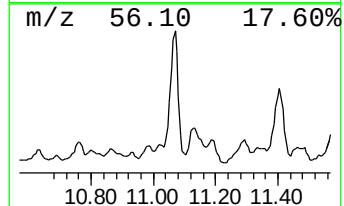
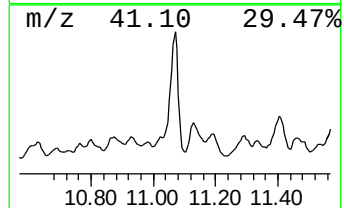
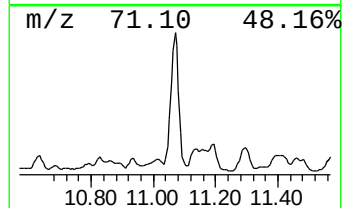
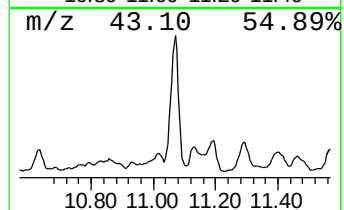
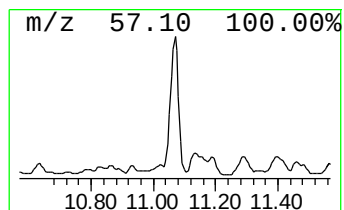
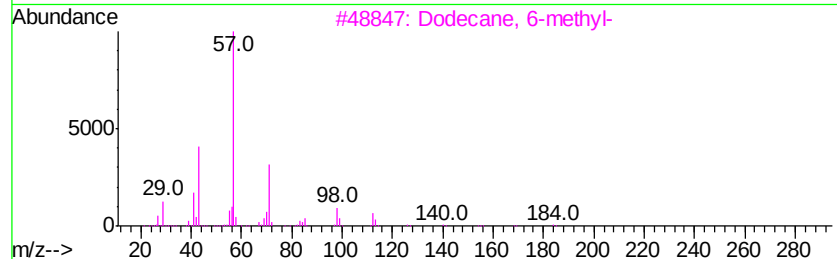
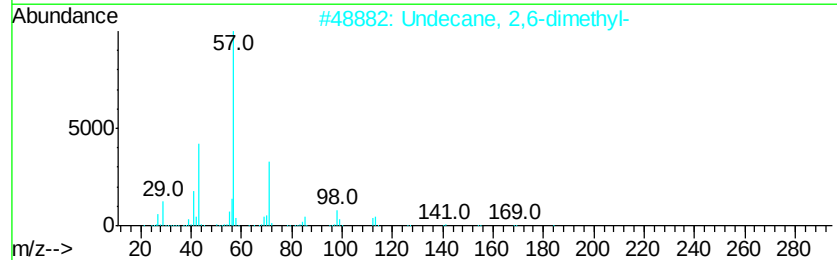
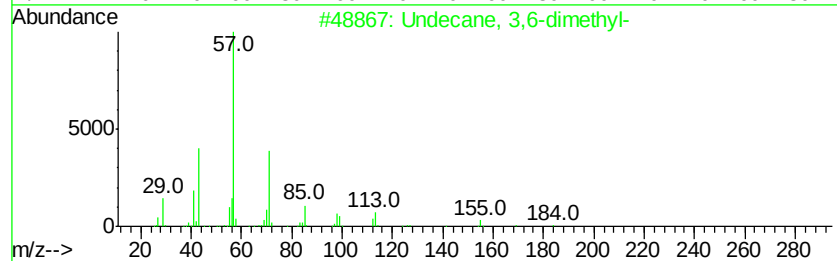
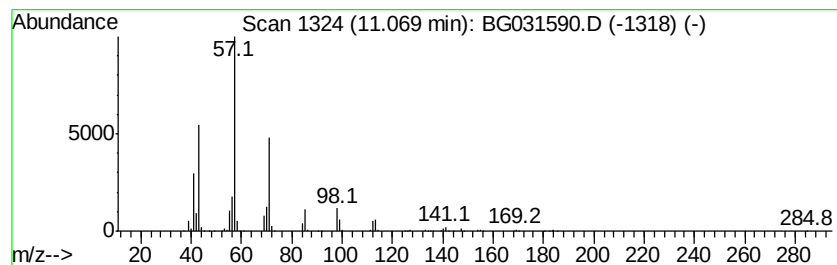
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Undecane, 3,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.07	89.66 ng	6698220	Napthalene-d8	10.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	91
2		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	90
3		Dodecane, 6-methyl-	184	C13H28	006044-71-9	76
4		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
5		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	64



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Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
Sample : I6830-04
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MAP-5-E(12-13)

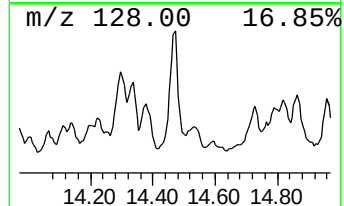
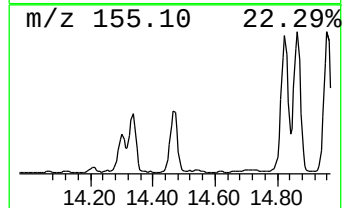
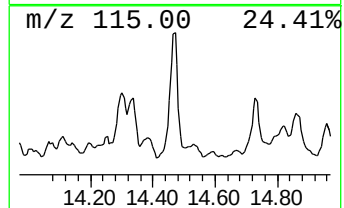
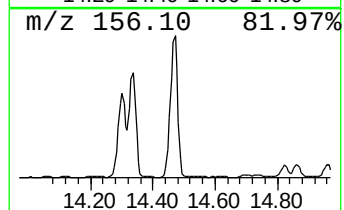
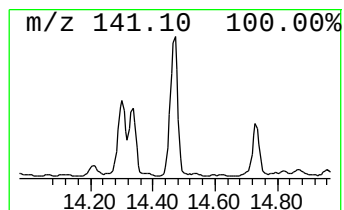
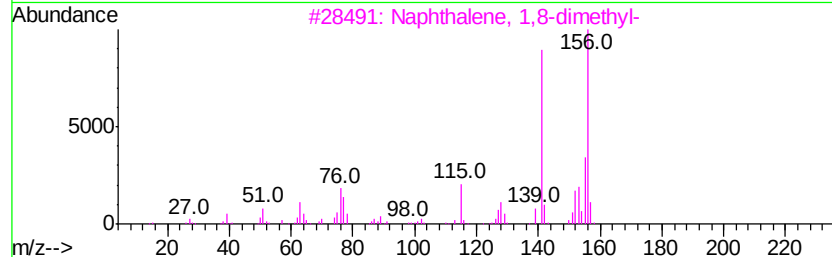
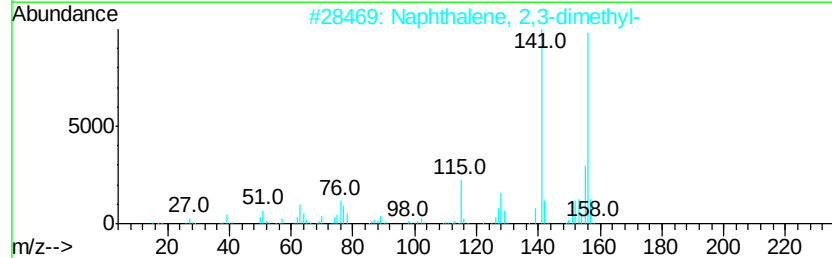
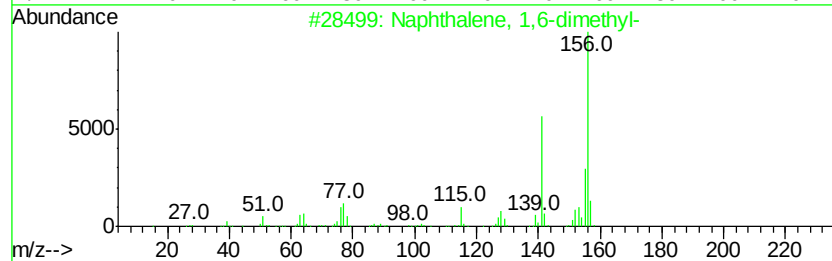
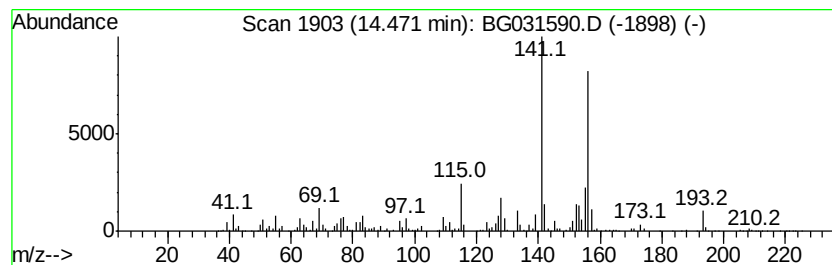
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	51.74 ng	7606810	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG121417\
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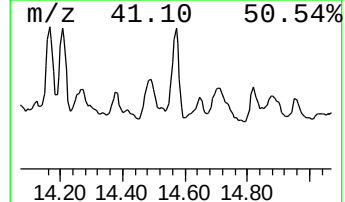
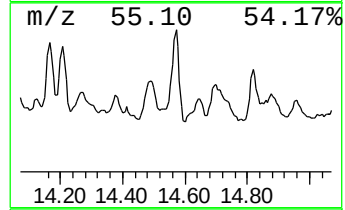
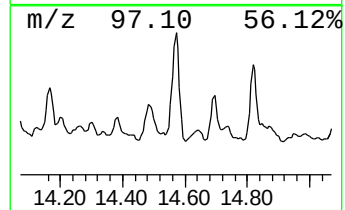
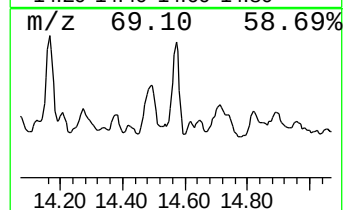
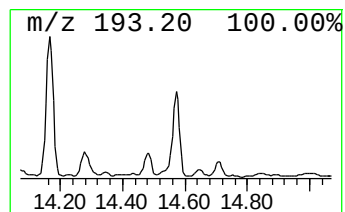
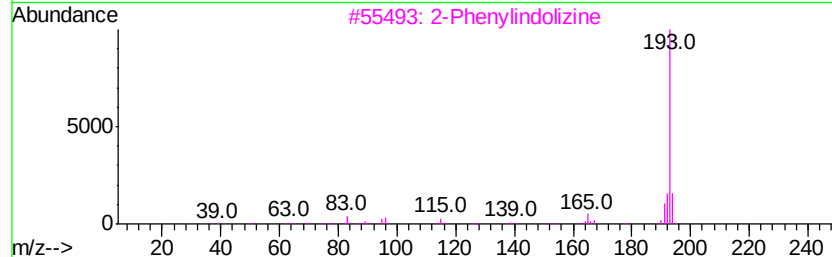
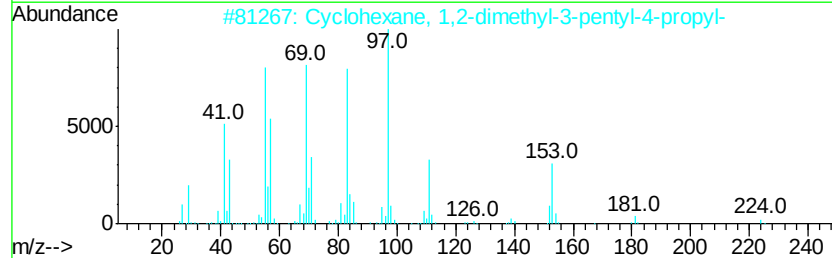
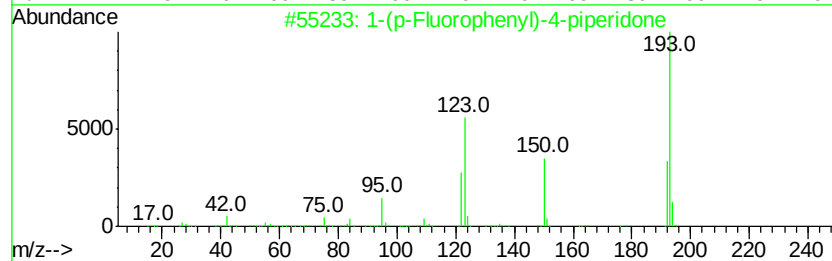
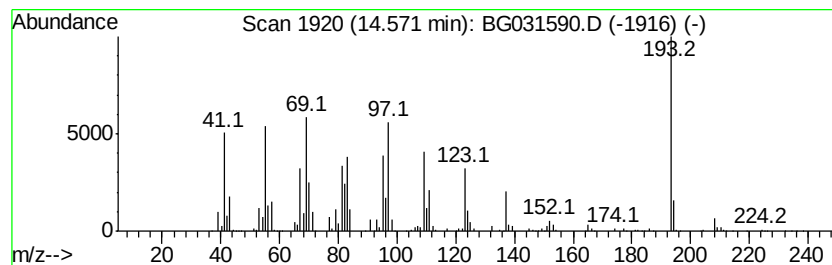
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-(p-Fluorophenyl)-4-piperi... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.57	44.40 ng	6527030	Acenaphthene-d10	14.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	35
2	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	35
3	2-Phenylindolizine	193	C14H11N	025379-20-8	35
4	Butan-2-one, 1-[3-(3,3-dimethyl-...	278	C16H26N2O2	1000303-34-2	22
5	6-Nitroundec-5-ene	199	C11H21NO2	1000192-40-3	22



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG121417\
Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
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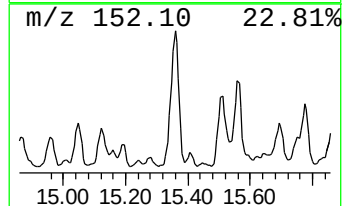
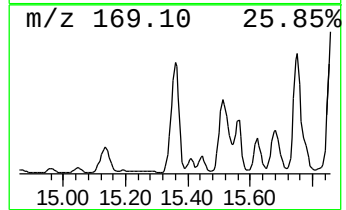
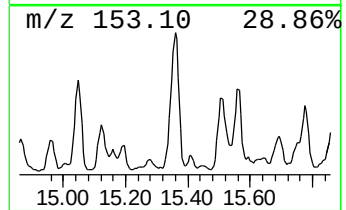
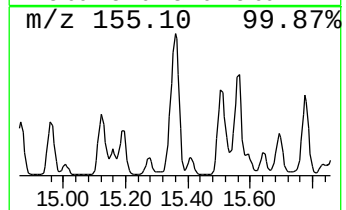
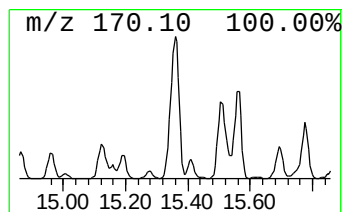
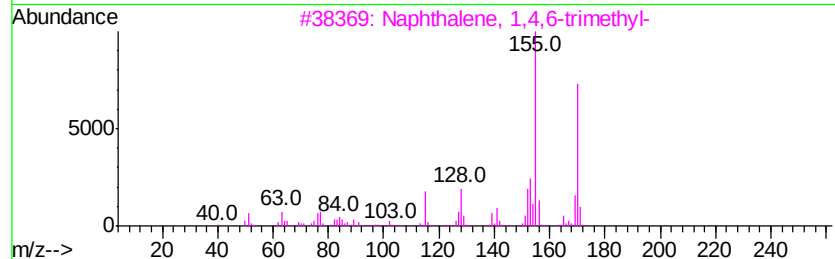
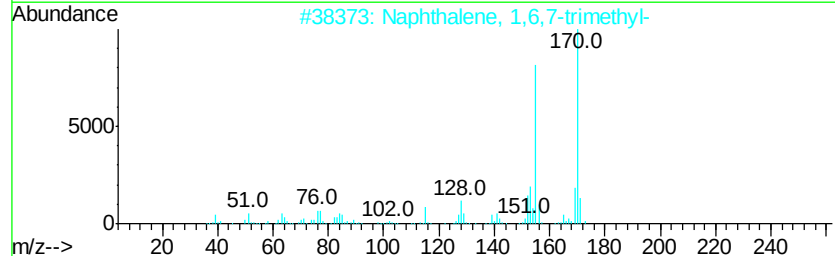
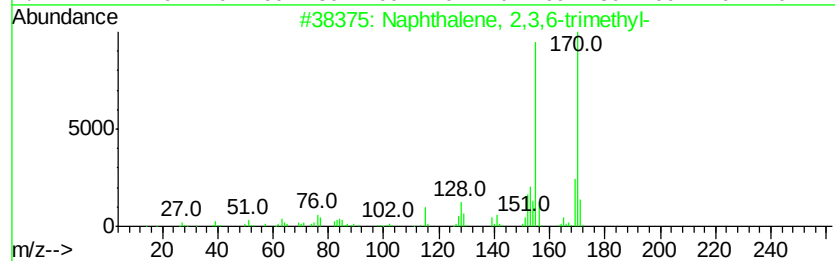
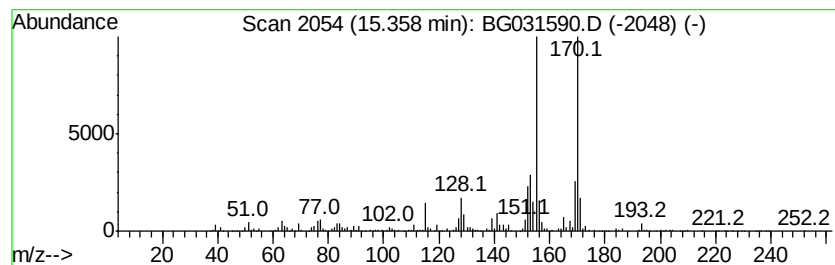
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 2,3,6-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	82.39 ng	12113200	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
4		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	93
5		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	93



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
Sample : I6830-04
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MAP-5-E(12-13)

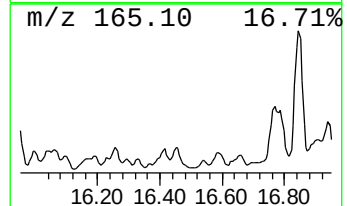
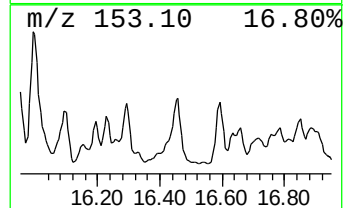
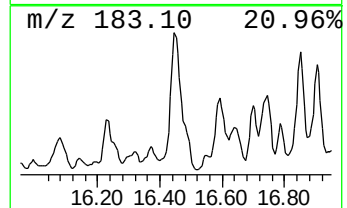
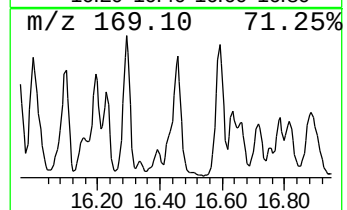
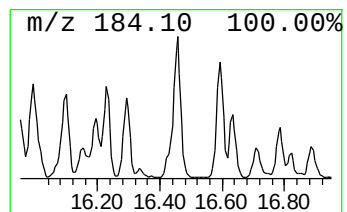
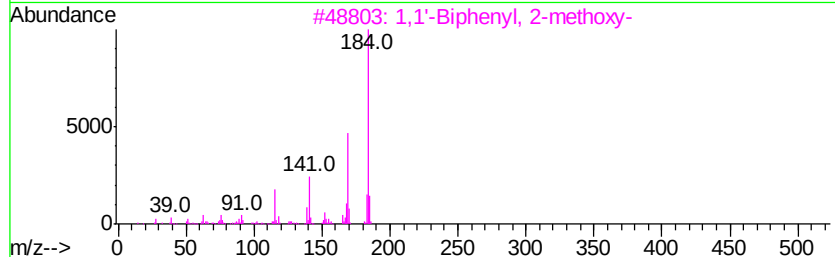
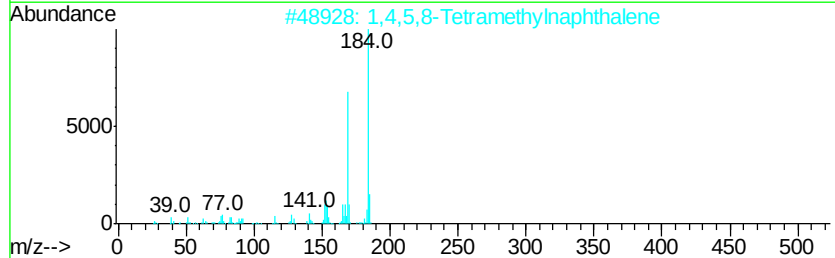
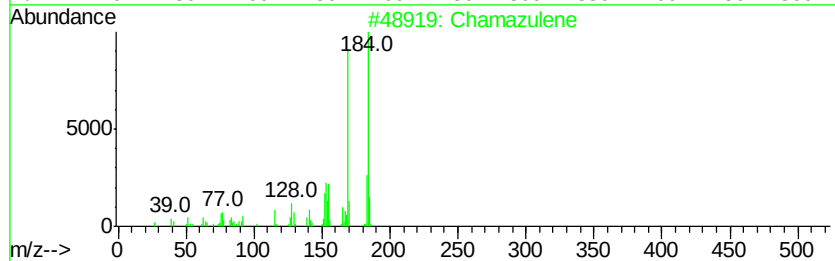
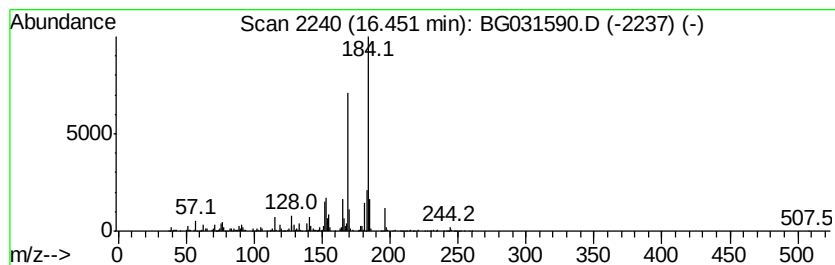
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Chamazulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.45	82.02 ng	6361990	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chamazulene	184	C14H16	000529-05-5	87
2		1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	81
3		1,1'-Biphenyl, 2-methoxy-	184	C13H12O	000086-26-0	64
4		Benzene, (4,5,5-trimethyl-1,3-cy...	184	C14H16	033930-85-7	64
5		Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	64



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG121417\
Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
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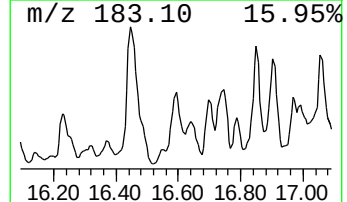
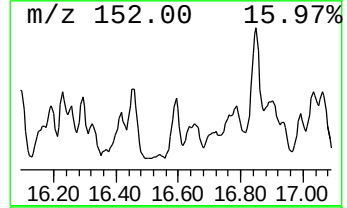
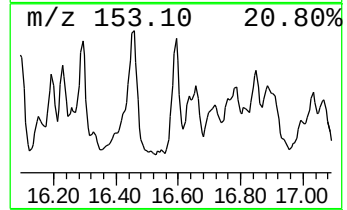
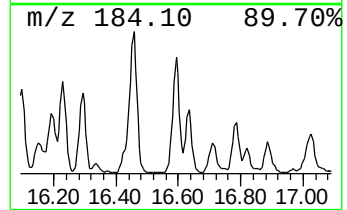
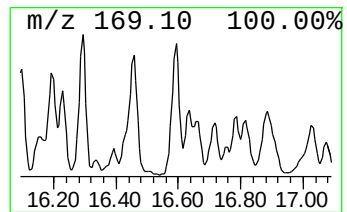
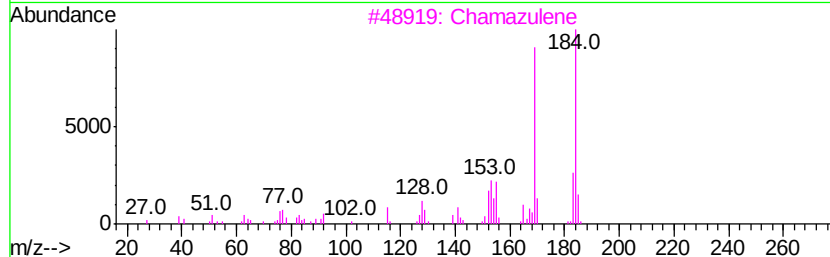
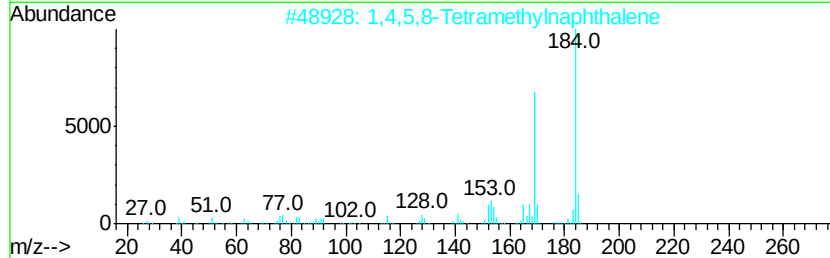
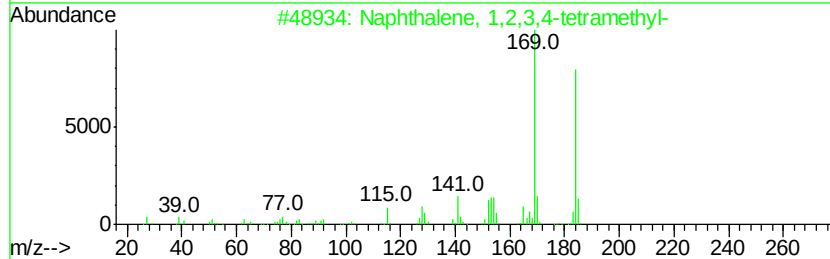
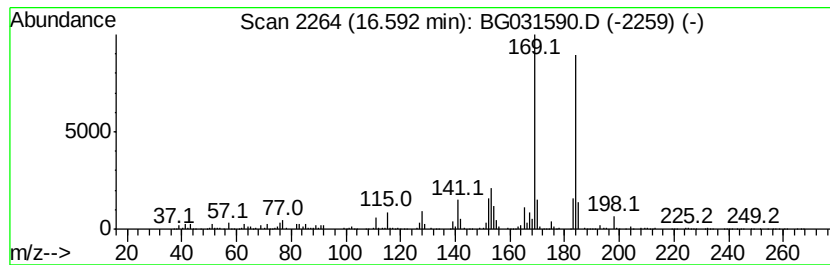
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 1,2,3,4-tetram... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	50.57 ng	3922040	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	93
2		1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	93
3		Chamazulene	184	C14H16	000529-05-5	93
4		Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	91
5		Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	86



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG121417\
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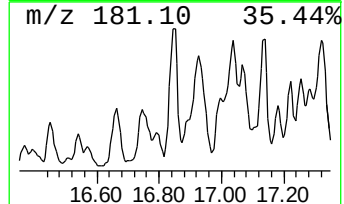
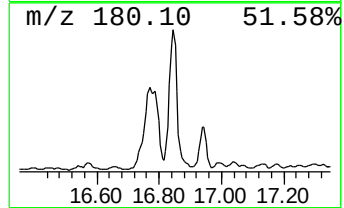
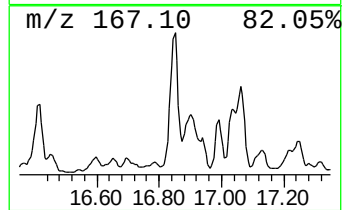
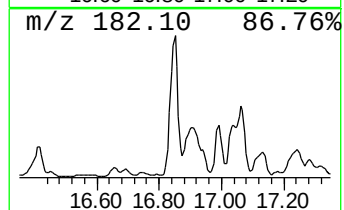
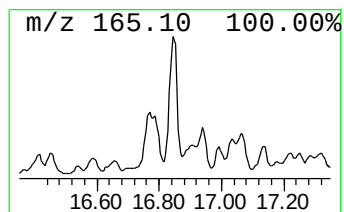
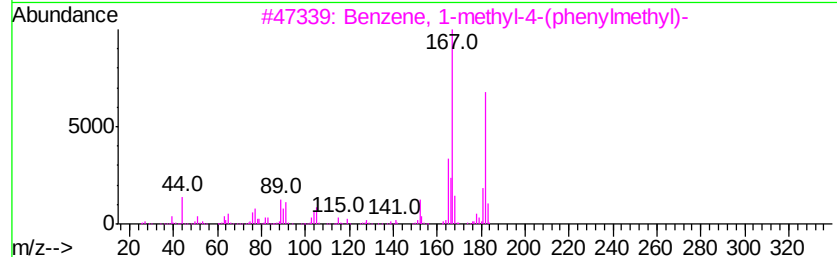
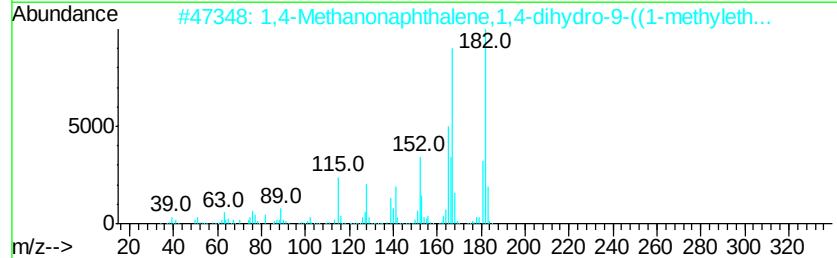
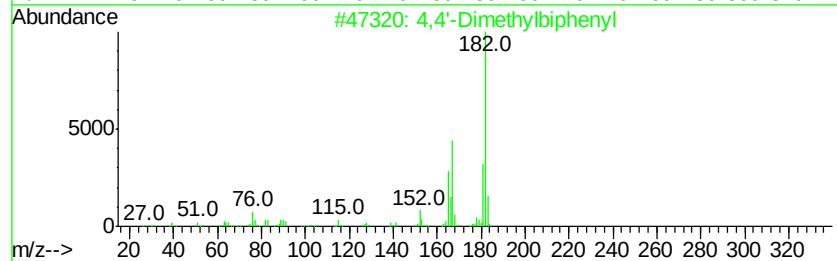
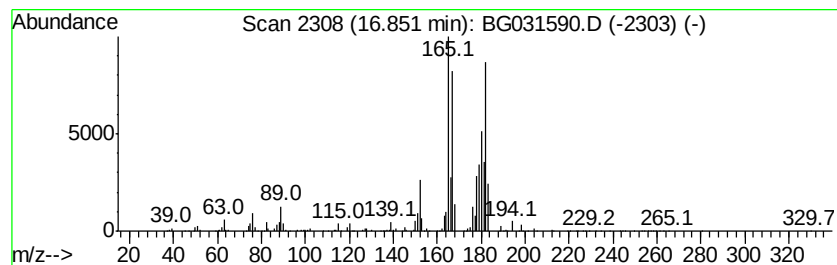
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 4,4'-Dimethylbiphenyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.85	71.91 ng	5577690	Phenanthrene-d10	17.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4,4'-Dimethylbiphenyl	182 C14H14	000613-33-2 62
2		1,4-Methanonaphthalene,1,4-dihyd...	182 C14H14	007350-72-3 58
3		Benzene, 1-methyl-4-(phenylmethyl)-	182 C14H14	000620-83-7 49
4		1,1'-Biphenyl, 2-ethyl-	182 C14H14	001812-51-7 49
5		3H-Benz[e]indene, 2-methyl-	180 C14H12	150096-60-9 42



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
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ALS Vial : 29 Sample Multiplier: 1

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ClientSampleId :
MAP-5-E(12-13)

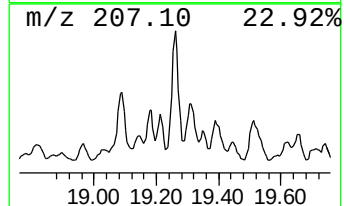
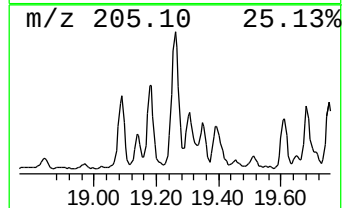
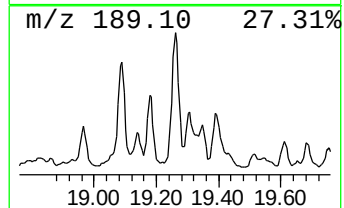
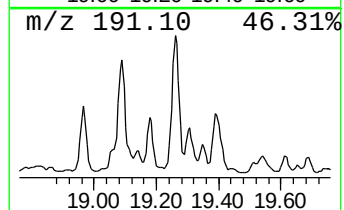
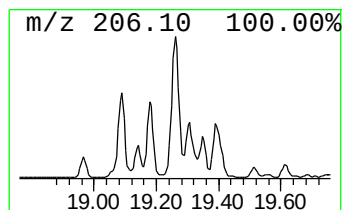
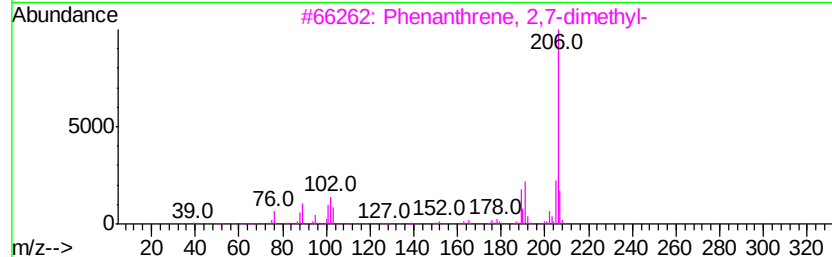
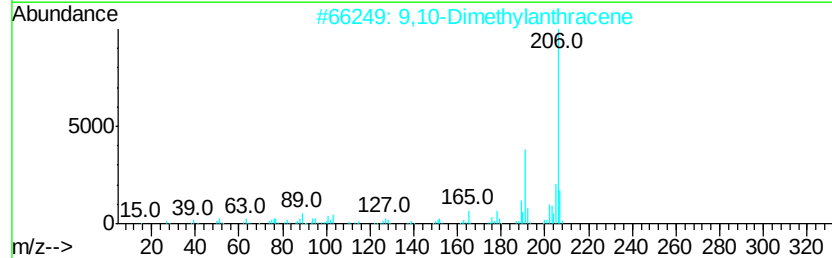
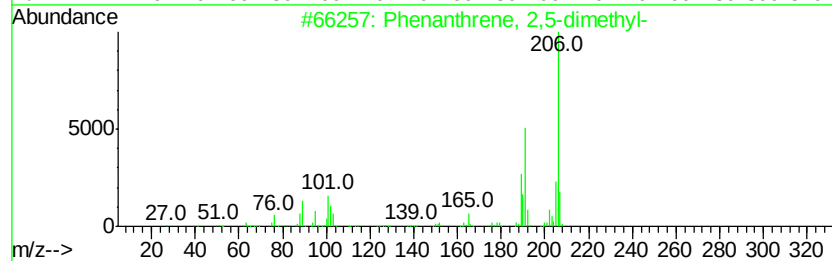
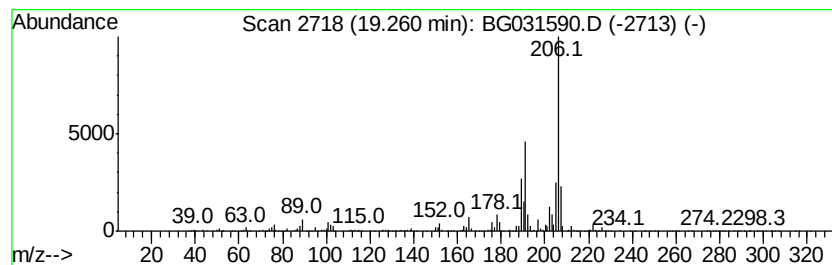
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	44.78 ng	3473230	Phenanthrene-d10	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	90
4		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG121417\
Data File : BG031590.D
Acq On : 15 Dec 2017 11:30
Operator : SJ/JU
Sample : I6830-04
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MAP-5-E(12-13)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 4-ethyl-...	9.20	59.0	ng	3415630	1	8.10	1157320	20.0
Naphthalene, deca...	10.09	50.8	ng	3794110	2	10.92	1494150	20.0
Undecane, 3,6-dim...	11.07	89.7	ng	6698220	2	10.92	1494150	20.0
Naphthalene, 1,6-...	14.47	51.7	ng	7606810	3	14.74	2940350	20.0
1-(p-Fluorophenyl...	14.57	44.4	ng	6527030	3	14.74	2940350	20.0
Naphthalene, 2,3,...	15.36	82.4	ng	12113200	3	14.74	2940350	20.0
Chamazulene	16.45	82.0	ng	6361990	4	17.49	1551240	20.0
Naphthalene, 1,2,...	16.59	50.6	ng	3922040	4	17.49	1551240	20.0
4,4'-Dimethylbiph...	16.85	71.9	ng	5577690	4	17.49	1551240	20.0
Phenanthrene, 2,5...	19.26	44.8	ng	3473230	4	17.49	1551240	20.0