

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
 Data File : BF127156.D
 Acq On : 02 Mar 2022 19:35
 Operator : CG\JU
 Sample : N1708-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PPSP-B10-17.5-18.0

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-BF021622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127156.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.516	545	549	552	rVB	2926592	3109481	17.10%	0.679%
2	6.134	649	654	660	rBV2	1552298	2595145	14.27%	0.566%
3	6.328	681	687	689	rBV3	927849	1331644	7.32%	0.291%
4	6.387	694	697	699	rVB	1588760	1387169	7.63%	0.303%
5	6.416	699	702	705	rBV2	1852595	2321939	12.77%	0.507%
6	6.528	715	721	725	rBV	2599231	3617041	19.89%	0.790%
7	6.634	734	739	743	rBV5	1130792	2186018	12.02%	0.477%
8	6.722	750	754	757	rBV	4408334	4643656	25.53%	1.014%
9	6.898	781	784	787	rVB2	2879966	3310065	18.20%	0.723%
10	6.945	787	792	795	rBV2	955488	1304646	7.17%	0.285%
11	7.040	804	808	811	rBV4	2162625	2591692	14.25%	0.566%
12	7.075	811	814	819	rVB3	1424606	1513664	8.32%	0.330%
13	7.140	823	825	827	rBV	1344934	1368538	7.53%	0.299%
14	7.163	827	829	832	rVB	2307399	2375403	13.06%	0.519%
15	7.216	832	838	840	rBV2	3152640	5070599	27.88%	1.107%
16	7.287	844	850	853	rBV3	3368272	5396203	29.67%	1.178%
17	7.363	859	863	865	rBV	2144423	2565760	14.11%	0.560%
18	7.387	865	867	870	rVB	1873643	1729364	9.51%	0.377%
19	7.440	870	876	878	rBV	5405207	7296751	40.12%	1.593%
20	7.469	878	881	884	rVB2	4489327	4358356	23.97%	0.951%
21	7.540	888	893	896	rBV4	3203978	4805107	26.42%	1.049%
22	7.604	896	904	908	rBV5	2319808	5478912	30.13%	1.196%
23	7.675	908	916	918	rBV3	3411941	5582752	30.70%	1.219%
24	7.698	918	920	924	rVB2	4431959	4494492	24.71%	0.981%
25	7.787	930	935	936	rBV2	2511697	3267699	17.97%	0.713%
26	7.804	936	938	940	rVV2	3161171	3454786	19.00%	0.754%
27	7.828	940	942	945	rVV3	3098195	4846567	26.65%	1.058%
28	7.857	945	947	951	rVV3	3997164	5467691	30.07%	1.194%
29	7.898	951	954	956	rVV3	2581142	3092297	17.00%	0.675%
30	7.928	956	959	962	rVV3	6223048	7756211	42.65%	1.693%
31	7.957	962	964	966	rVV	3510941	3185610	17.52%	0.695%
32	7.981	966	968	969	rVV	2184031	1913030	10.52%	0.418%
33	8.004	969	972	974	rVV2	3414210	4162675	22.89%	0.909%
34	8.034	974	977	979	rVV	3787980	3404808	18.72%	0.743%
35	8.057	979	981	984	rVB2	2761265	2357982	12.97%	0.515%
36	8.145	990	996	998	rBV7	2788266	5169547	28.43%	1.128%

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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-BF021622.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	8.181	999	1002	1004	rVV	5179876	6683887	36.75%	1.459%
38	8.216	1004	1008	1010	rVV3	6325802	9385980	51.61%	2.049%
39	8.239	1010	1012	1016	rVV2	5486925	7524094	41.37%	1.642%
40	8.275	1016	1018	1021	rVB	2475842	2132674	11.73%	0.466%
41	8.345	1026	1030	1031	rBV4	1711088	2130785	11.72%	0.465%
42	8.392	1035	1038	1041	rVB2	5764695	5707225	31.38%	1.246%
43	8.439	1042	1046	1048	rBV	3525782	3465904	19.06%	0.757%
44	8.481	1048	1053	1058	rBV5	3932428	7077567	38.92%	1.545%
45	8.539	1060	1063	1066	rVB4	2333536	2305314	12.68%	0.503%
46	8.581	1066	1070	1072	rBV2	3835104	3933633	21.63%	0.859%
47	8.622	1072	1077	1079	rBV2	4238430	4881478	26.84%	1.066%
48	8.710	1087	1092	1095	rBV	7101029	9409818	51.74%	2.054%
49	8.792	1103	1106	1108	rBV2	1899784	2423914	13.33%	0.529%
50	8.863	1116	1118	1124	rVB4	3014237	4579256	25.18%	1.000%
51	8.934	1124	1130	1132	rVB	6585370	11316946	62.23%	2.470%
52	8.975	1135	1137	1140	rBV3	979101	1231141	6.77%	0.269%
53	9.034	1140	1147	1151	rBV4	8925213	16381342	90.08%	3.576%
54	9.145	1162	1166	1168	rBV4	3250430	3383805	18.61%	0.739%
55	9.169	1168	1170	1173	rVB2	2764650	2199902	12.10%	0.480%
56	9.222	1175	1179	1181	rVV3	4940691	6396906	35.18%	1.396%
57	9.275	1185	1188	1190	rBV2	2649163	2311976	12.71%	0.505%
58	9.351	1198	1201	1203	rBV4	2393501	3125326	17.19%	0.682%
59	9.404	1208	1210	1211	rVV2	1744184	1605413	8.83%	0.350%
60	9.439	1211	1216	1218	rVV2	4100146	7089429	38.98%	1.548%
61	9.481	1218	1223	1229	rVV2	4234136	8776525	48.26%	1.916%
62	9.563	1230	1237	1240	rVV2	7000378	14006994	77.02%	3.057%
63	9.639	1240	1250	1253	rVV	6751654	18185538	100.00%	3.970%
64	9.675	1254	1256	1260	rVV4	8826429	12341692	67.87%	2.694%
65	9.722	1260	1264	1266	rVV3	2172879	3258597	17.92%	0.711%
66	9.751	1266	1269	1276	rVB4	6549010	11648397	64.05%	2.543%
67	9.834	1279	1283	1285	rVV2	4020686	4050813	22.27%	0.884%
68	9.875	1288	1290	1292	rVB2	2065258	1540645	8.47%	0.336%
69	9.945	1299	1302	1307	rVB3	4359625	5494944	30.22%	1.199%
70	10.004	1307	1312	1314	rBV3	3125526	5172866	28.44%	1.129%
71	10.075	1319	1324	1328	rBV2	4709433	8180813	44.99%	1.786%
72	10.110	1328	1330	1334	rBV3	1801984	1992525	10.96%	0.435%
73	10.175	1334	1341	1344	rVB2	5188054	7902437	43.45%	1.725%
74	10.216	1344	1348	1351	rBV3	5425710	6249841	34.37%	1.364%
75	10.286	1356	1360	1363	rBV2	5135402	7589965	41.74%	1.657%
76	10.316	1363	1365	1370	rVB3	4791016	5147277	28.30%	1.124%

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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-BF021622.M
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77	10.375	1370	1375	1377	rBV	3901796	5574160	30.65%	1.217%
78	10.398	1377	1379	1384	rVB4	2888328	3175962	17.46%	0.693%
79	10.463	1388	1390	1392	rVB	2126510	1662951	9.14%	0.363%
80	10.510	1395	1398	1400	rBV2	2243567	2441934	13.43%	0.533%
81	10.604	1409	1414	1416	rVV2	4729066	6385333	35.11%	1.394%
82	10.657	1420	1423	1425	rBV3	2147879	2623251	14.42%	0.573%
83	10.751	1437	1439	1443	rBV4	2125858	2707299	14.89%	0.591%
84	10.786	1443	1445	1447	rVB	1941964	1383202	7.61%	0.302%
85	10.875	1454	1460	1466	rVB2	8163097	12162338	66.88%	2.655%
86	10.945	1470	1472	1475	rVB2	1544904	1567566	8.62%	0.342%
87	11.045	1487	1489	1493	rBV3	1003714	1537278	8.45%	0.336%
88	11.098	1495	1498	1500	rVB2	2970638	2573228	14.15%	0.562%
89	11.175	1508	1511	1513	rBV2	1507830	1581509	8.70%	0.345%
90	11.333	1533	1538	1544	rVB2	5616506	8447400	46.45%	1.844%
91	11.480	1559	1563	1565	rBV	3008595	2692103	14.80%	0.588%
92	11.563	1570	1577	1579	rBV3	2095557	3498406	19.24%	0.764%
93	11.675	1593	1596	1607	rVB7	2535002	4456818	24.51%	0.973%
94	11.780	1609	1614	1619	rVV3	1487820	2822812	15.52%	0.616%
95	11.951	1640	1643	1645	rBV	2467287	2618797	14.40%	0.572%
96	11.986	1645	1649	1651	rVB	3360146	3592372	19.75%	0.784%
97	12.063	1659	1662	1664	rVV2	1608296	1632362	8.98%	0.356%
98	12.086	1664	1666	1671	rVB2	1835316	1911161	10.51%	0.417%
99	12.498	1732	1736	1739	rBV	2355862	2407964	13.24%	0.526%
100	13.022	1822	1825	1827	rBV	2226589	1927301	10.60%	0.421%

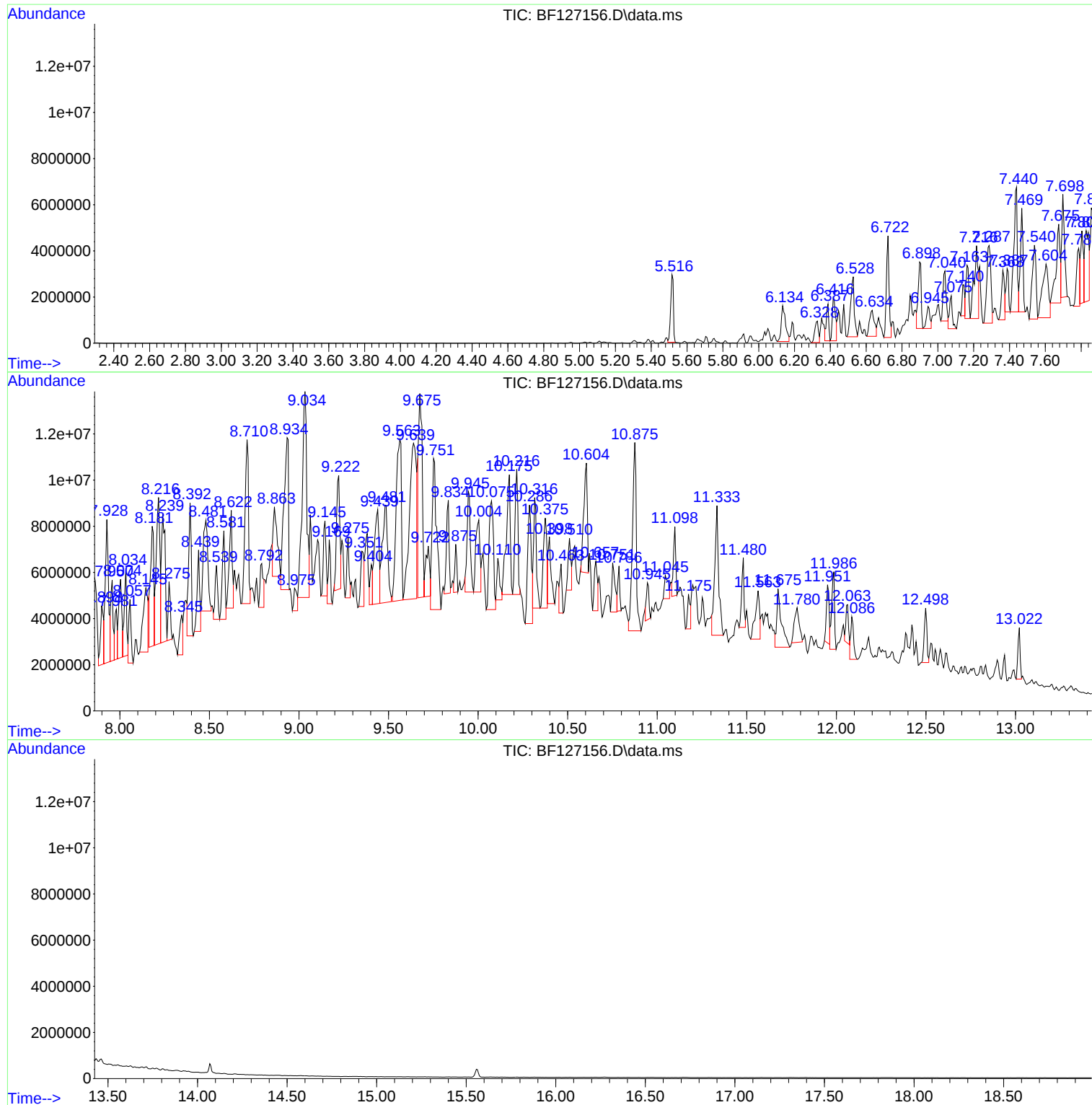
Sum of corrected areas: 458120391

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.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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PPSP-B10-17.5-18.0

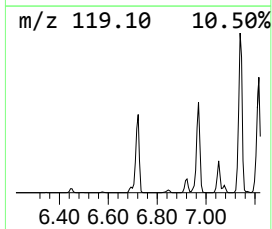
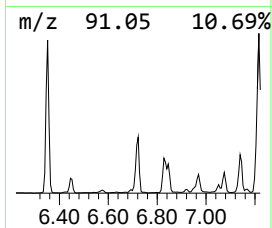
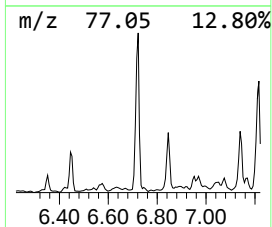
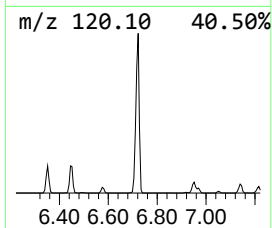
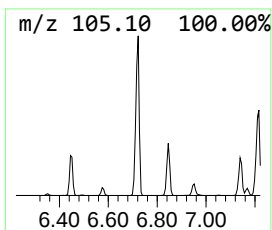
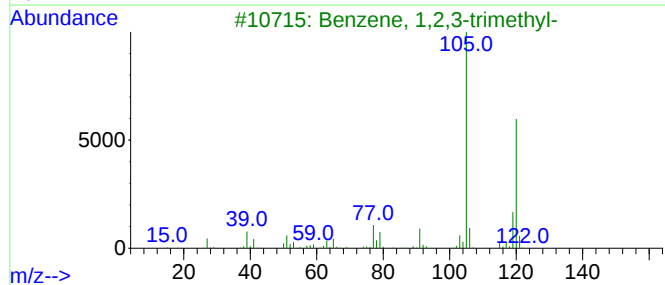
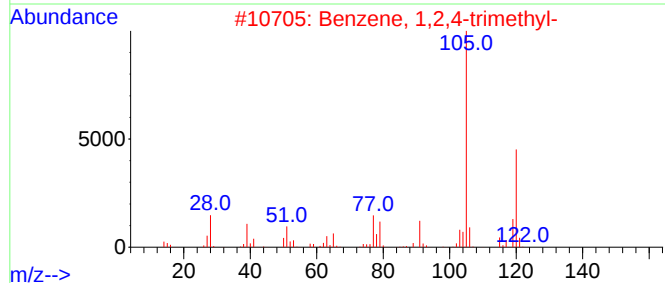
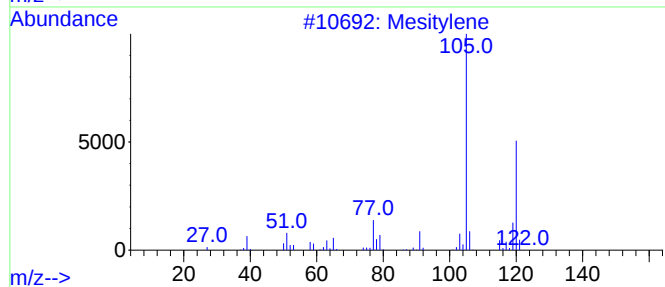
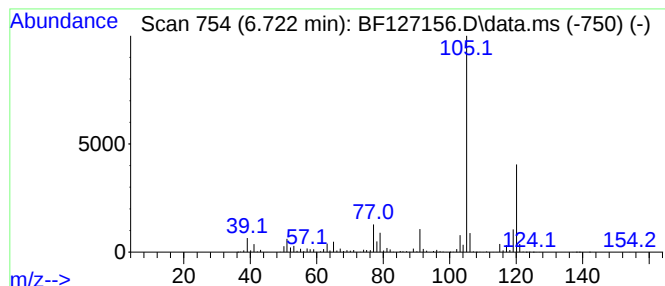
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Mesitylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.722	28.06 ng	4643660	1,4-Dichlorobenzene-d4	6.893

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Mesitylene	120	C9H12	000108-67-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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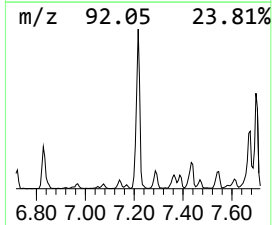
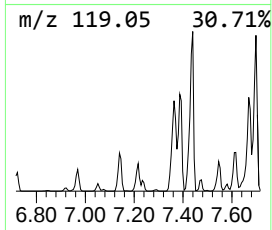
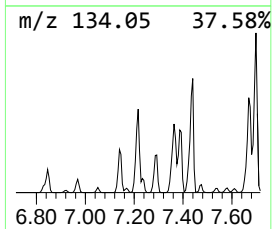
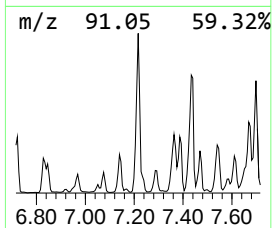
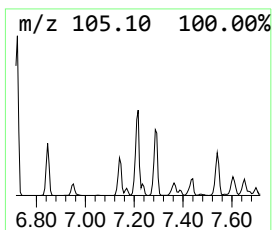
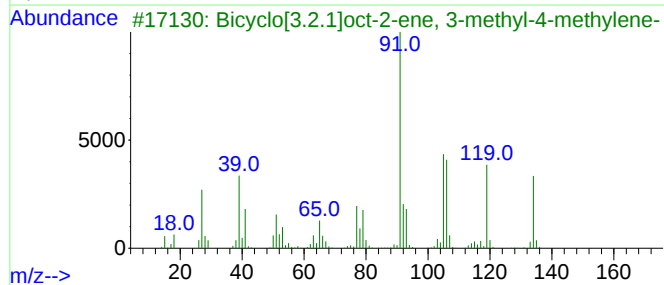
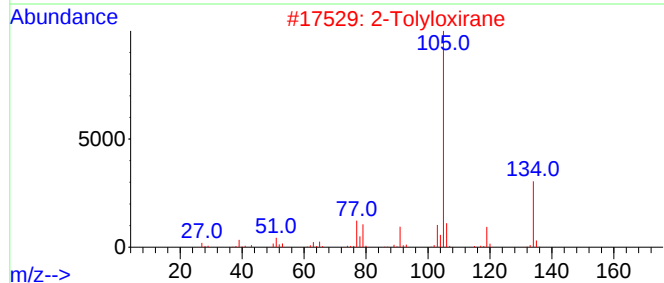
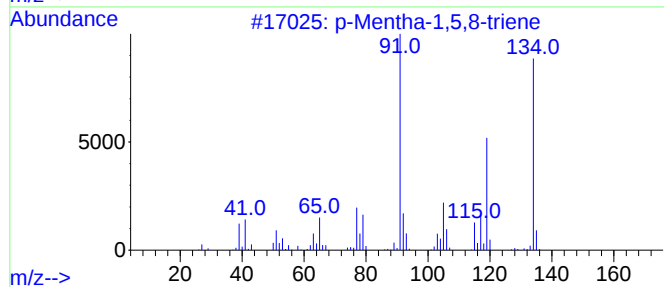
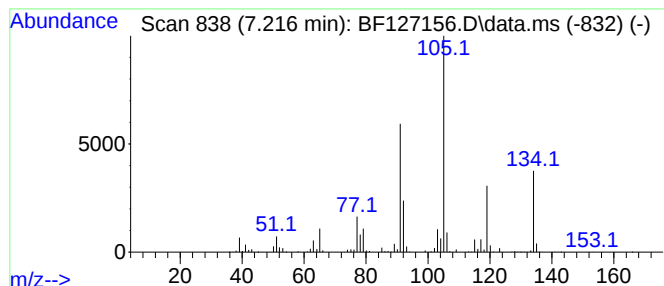
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 p-Mentha-1,5,8-triene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.216	30.64 ng	5070600	1,4-Dichlorobenzene-d4	6.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Mentha-1,5,8-triene	134	C10H14	021195-59-5	72
2			2-Tolyloxirane	134	C9H10O	002783-26-8	60
3			Bicyclo[3.2.1]oct-2-ene, 3-methy...	134	C10H14	049826-53-1	53
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	50
5			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	46



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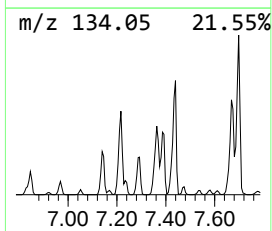
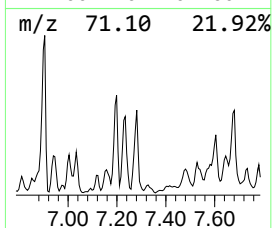
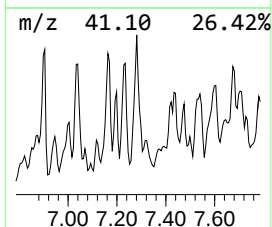
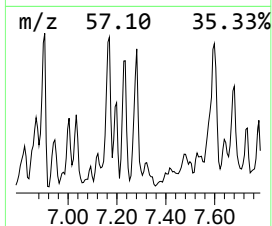
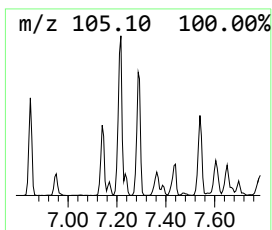
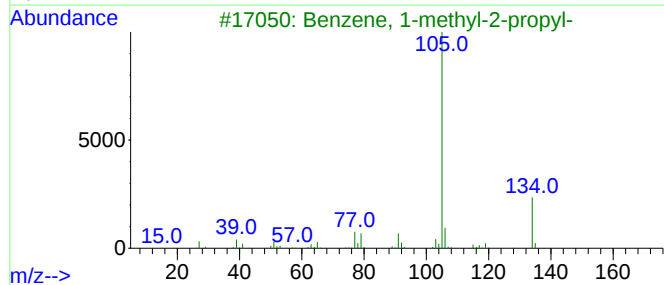
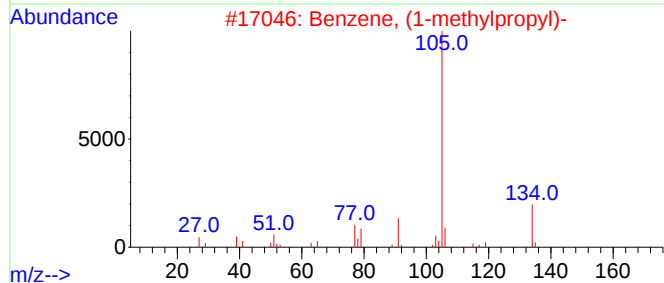
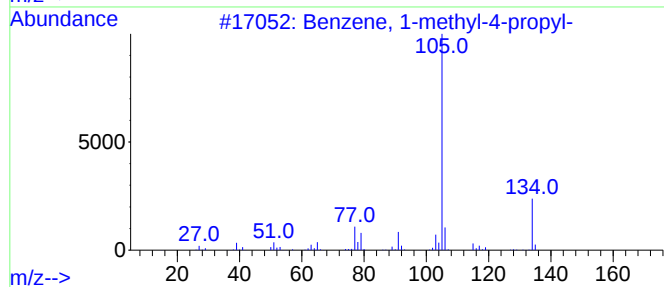
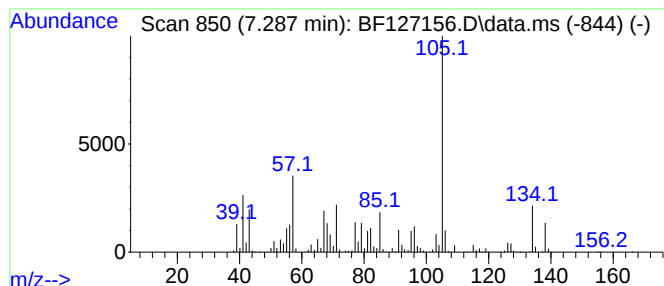
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, 1-methyl-4-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.287	32.60 ng	5396200	1,4-Dichlorobenzene-d4	6.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	74
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	70
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	38
4			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	38
5			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

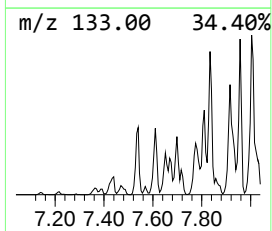
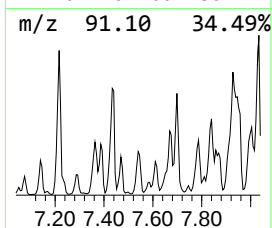
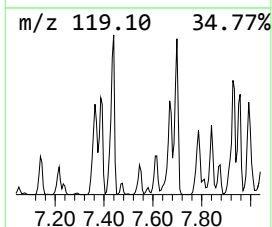
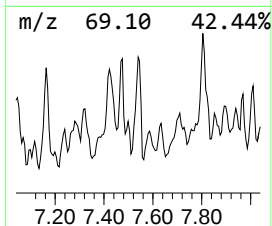
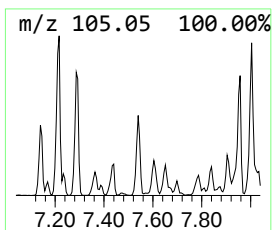
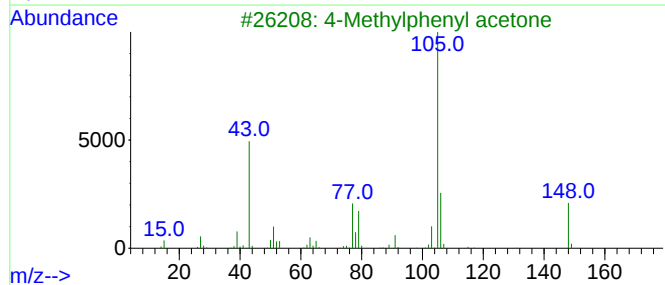
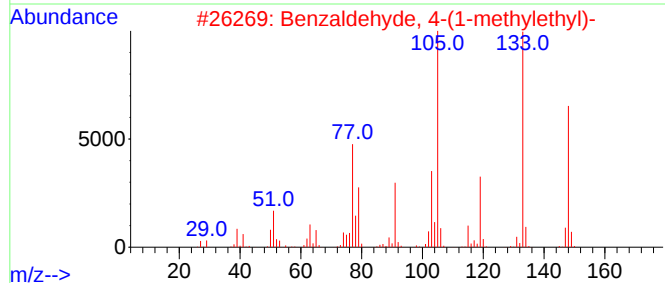
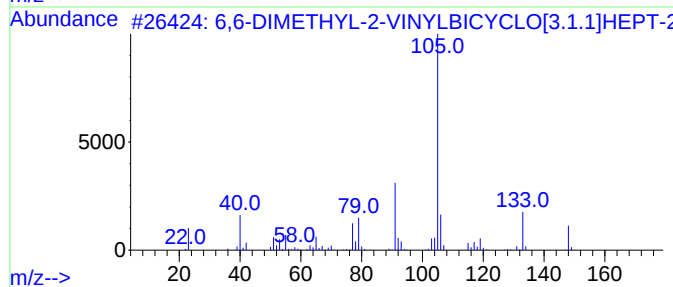
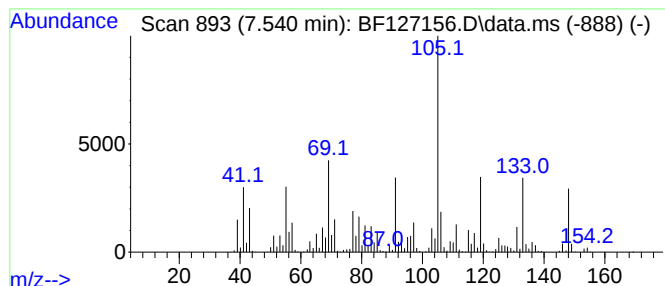
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ClientSampleId :
PPSP-B10-17.5-18.0

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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 6,6-DIMETHYL-2-VINYLBICYCLO... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.540	29.03 ng	4805110	1,4-Dichlorobenzene-d4	6.893	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6,6-DIMETHYL-2-VINYLBICYCLO[3.1....	148	C11H16	000473-00-7	53
2	Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	49
3	4-Methylphenyl acetone	148	C10H12O	002096-86-8	42
4	Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	42
5	Benzene, 1-methyl-4-butyl	148	C11H16	001595-05-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

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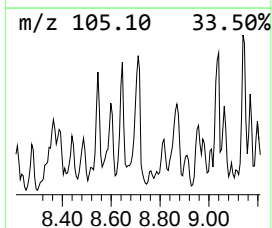
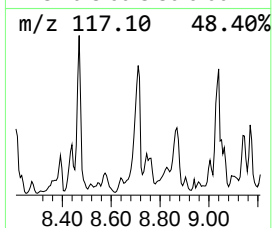
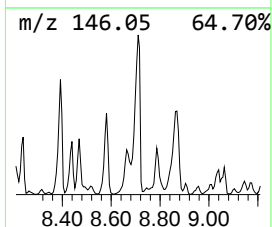
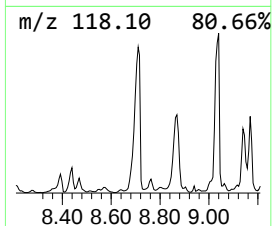
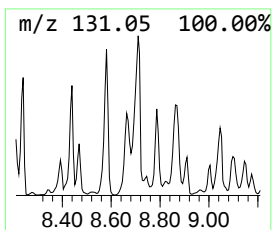
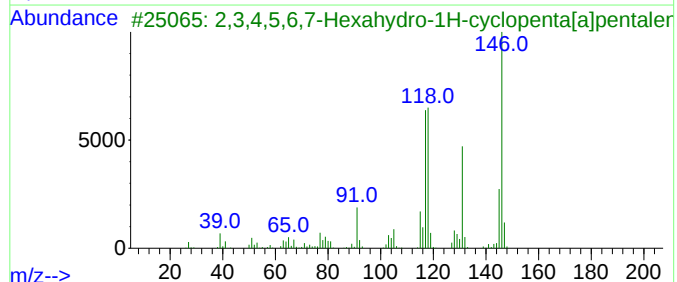
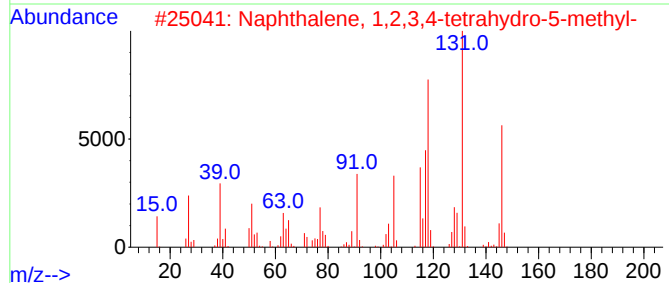
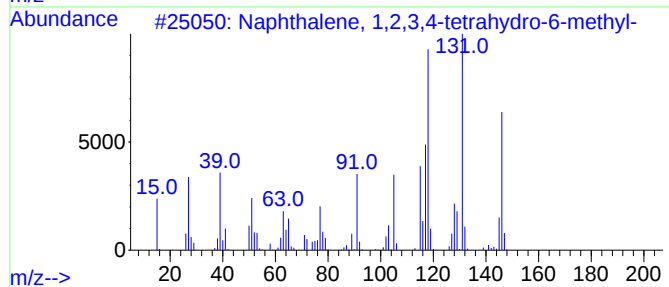
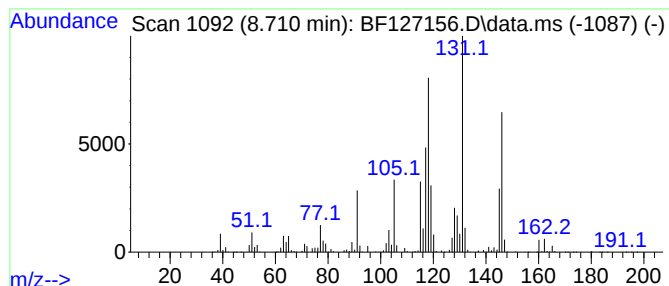
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.710	28.16 ng	9409820	Naphthalene-d8	8.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
3			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	90
4			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	72
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

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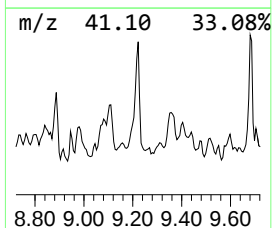
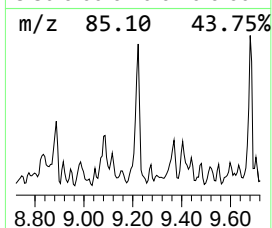
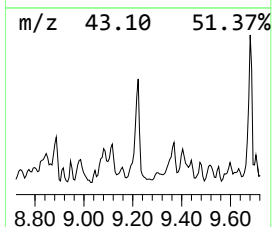
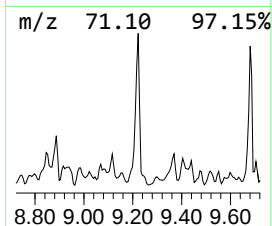
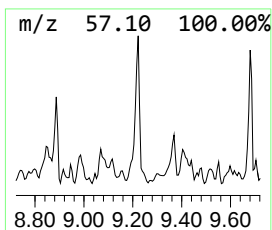
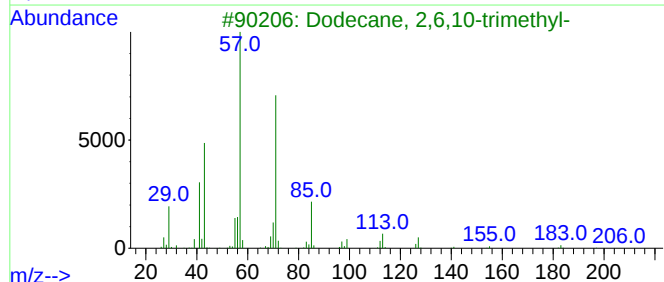
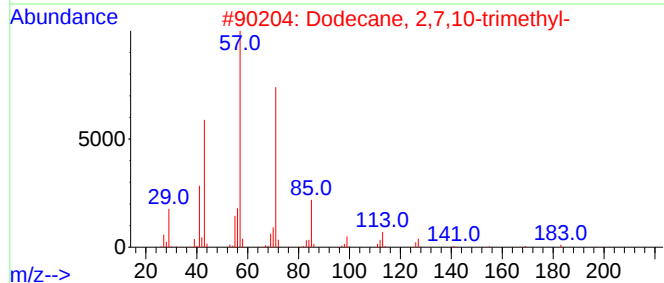
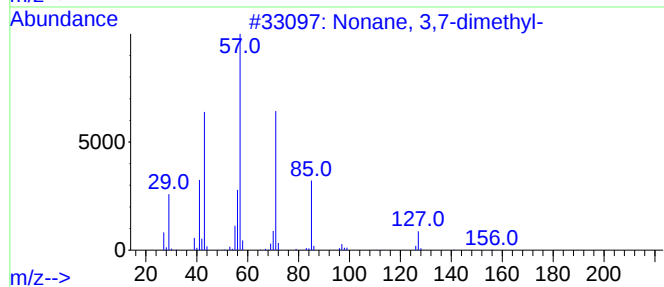
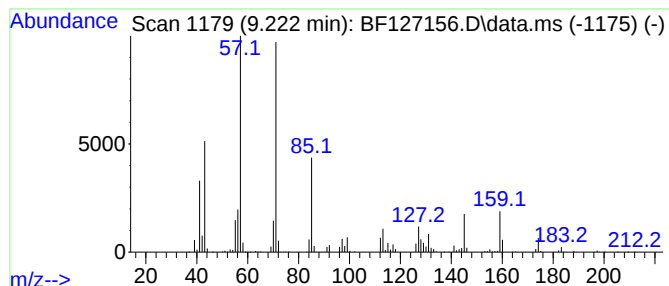
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Nonane, 3,7-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.222	23.28 ng	6396910	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	70
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	58
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	58
4			Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	49
5			Sulfurous acid, nonyl pentyl ester	278	C14H30O3S	1000309-14-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

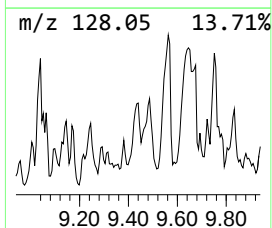
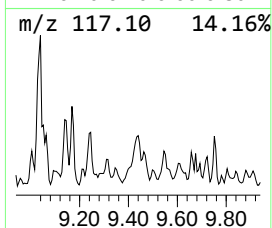
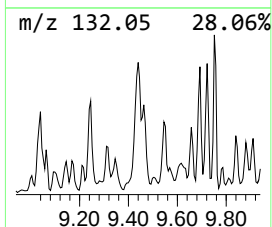
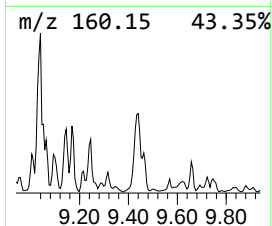
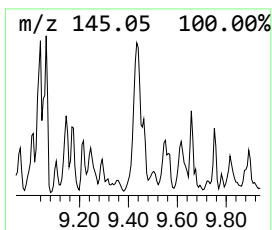
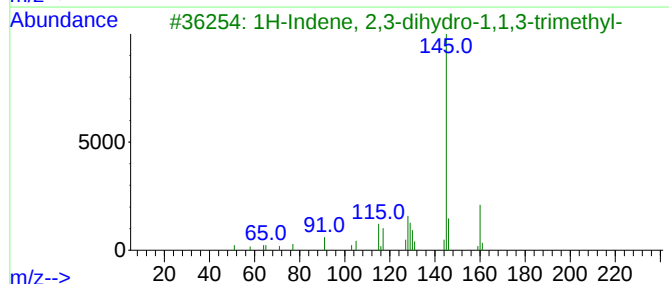
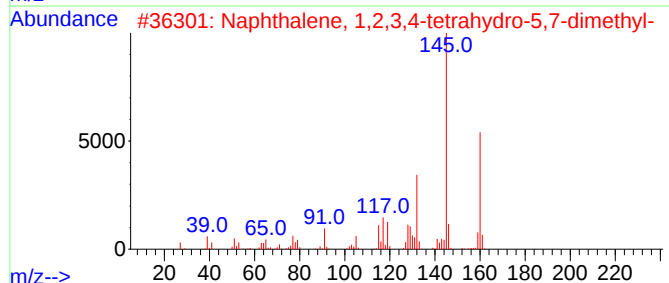
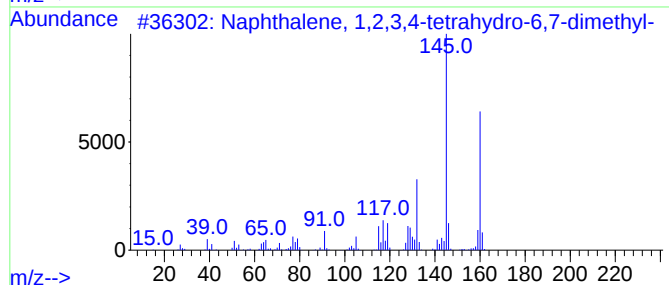
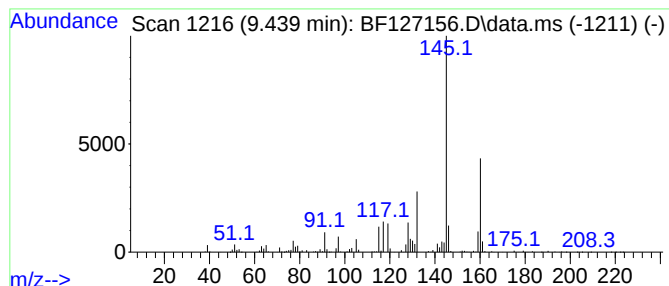
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.439	25.80 ng	7089430	Acenaphthene-d10	9.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	94
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	94
4	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	87
5	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

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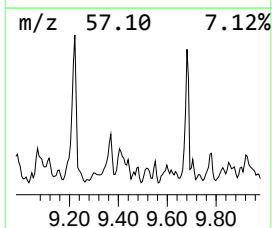
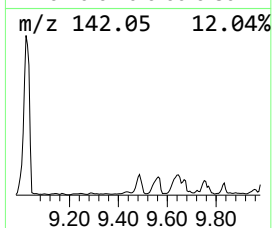
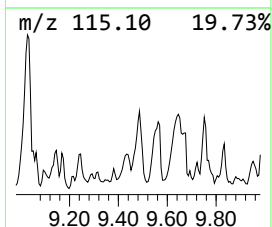
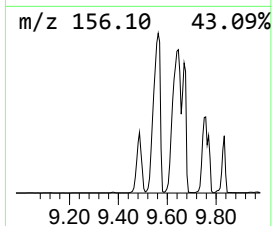
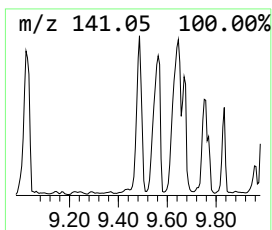
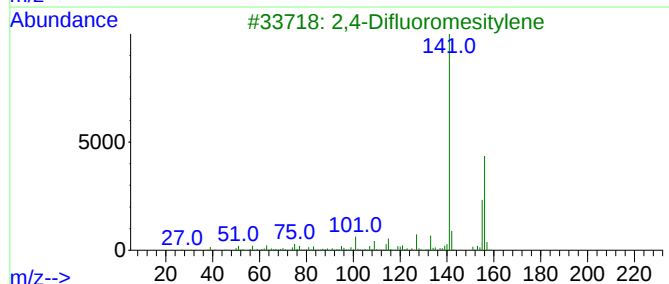
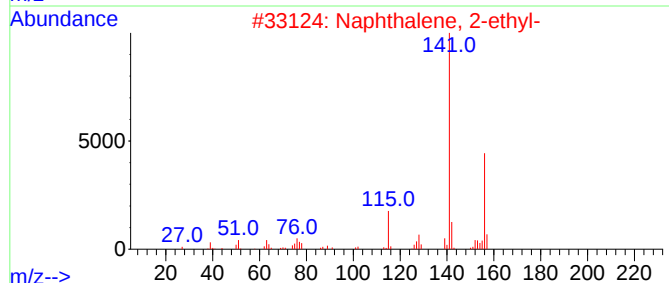
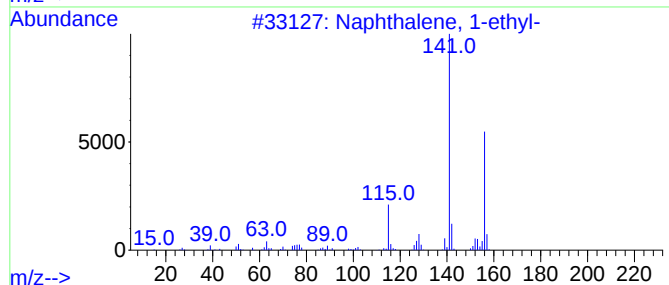
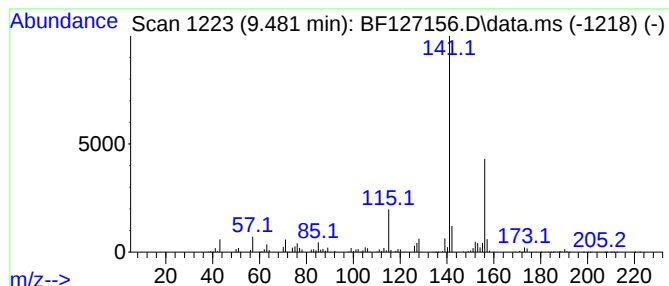
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.481	31.94 ng	8776530	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
2		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3		2,4-Difluoromesitylene	156	C9H10F2	000392-61-0	64
4		Butethal	212	C10H16N2O3	000077-28-1	59
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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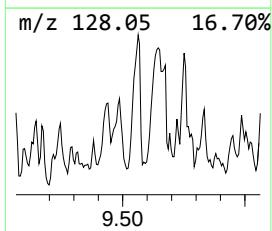
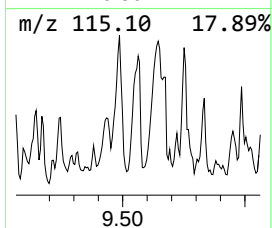
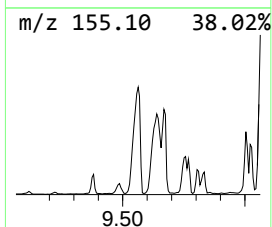
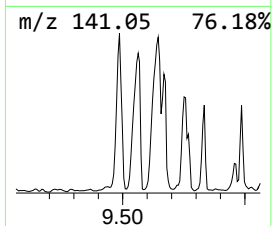
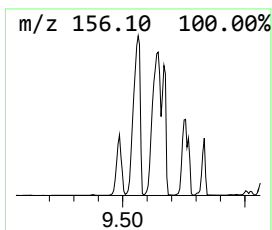
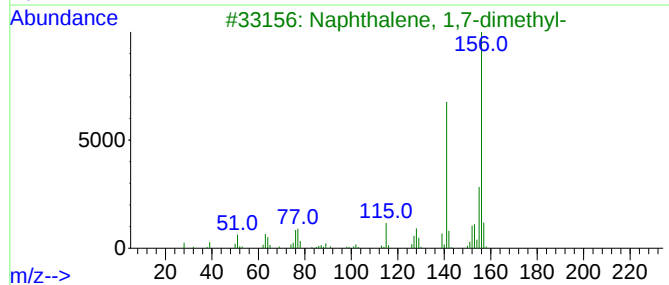
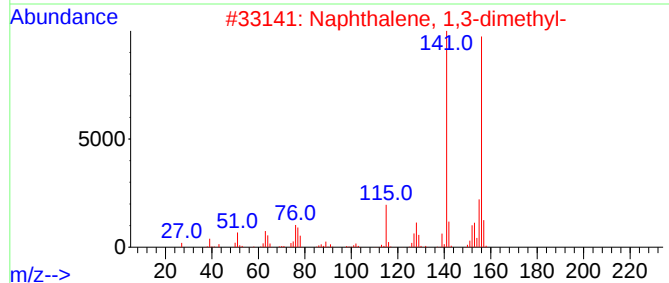
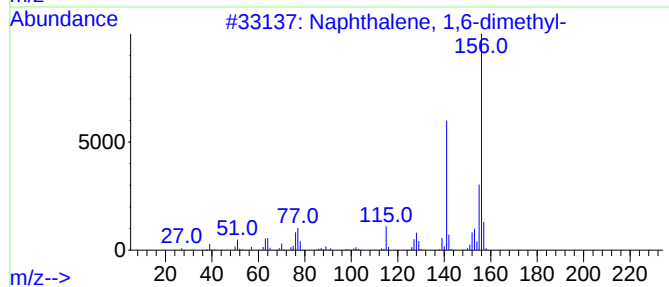
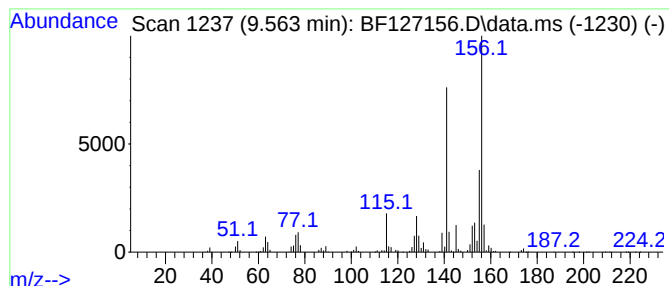
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.563	50.98 ng	14007000	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PPSP-B10-17.5-18.0

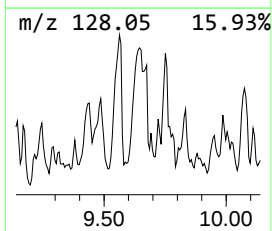
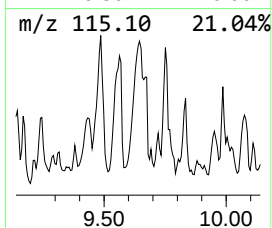
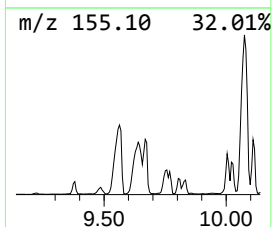
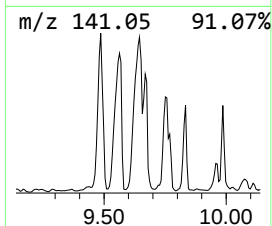
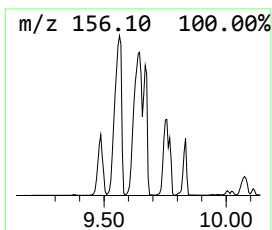
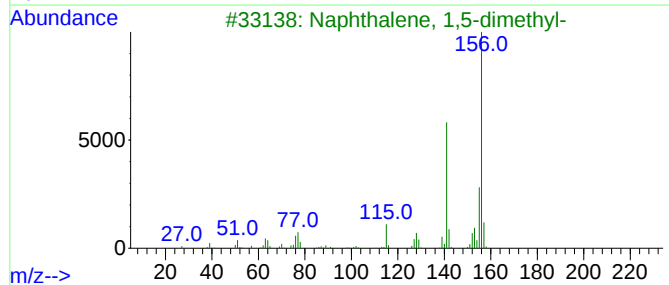
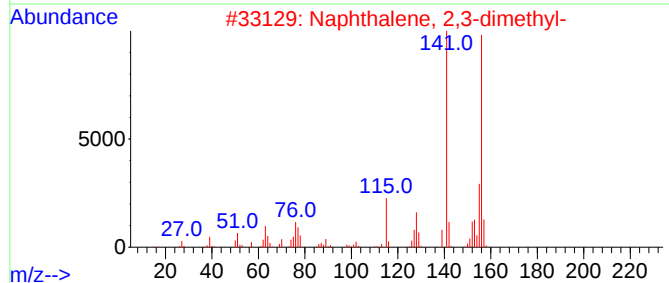
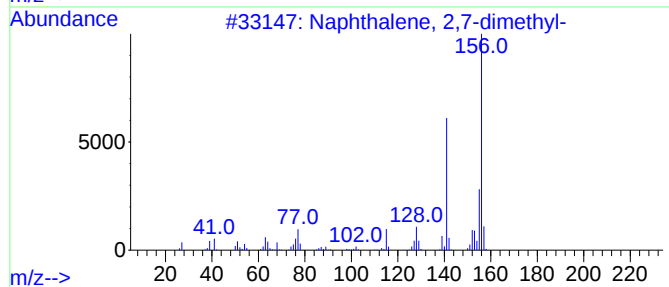
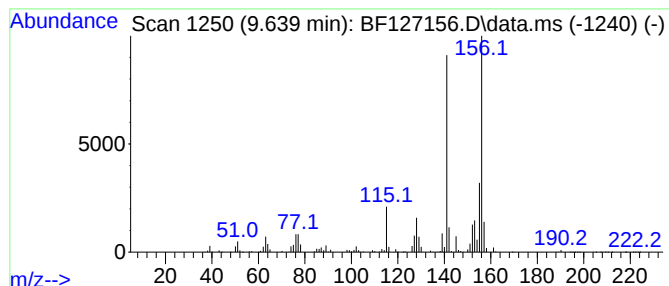
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.639	66.19 ng	18185500	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PPSP-B10-17.5-18.0

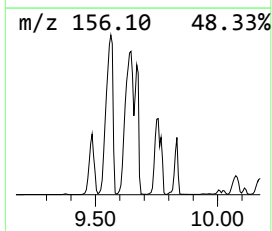
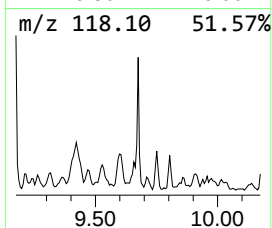
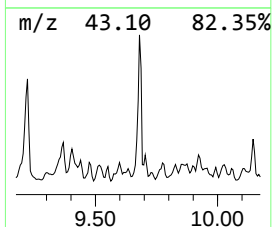
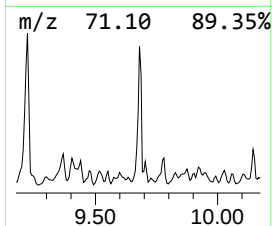
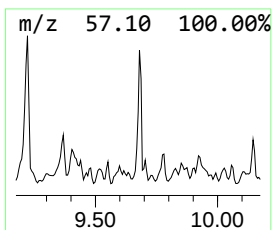
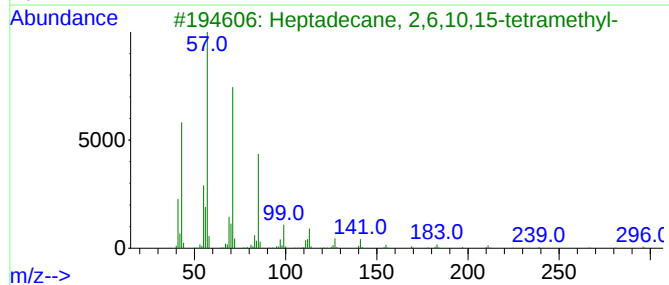
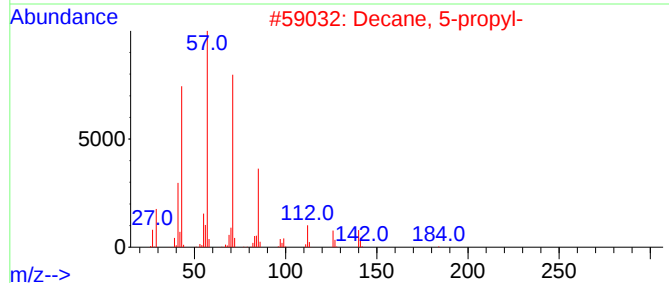
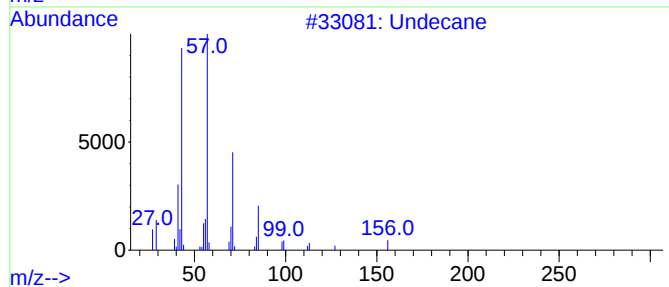
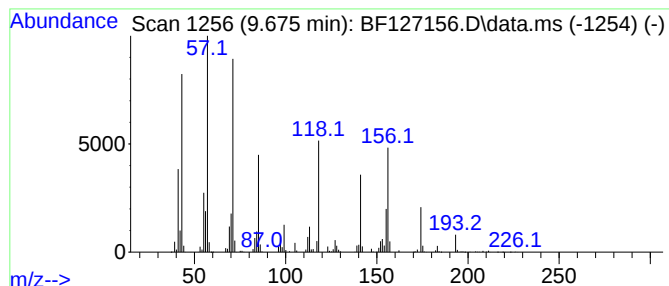
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Undecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.675	44.92 ng	12341700	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	60
2	Decane, 5-propyl-			184	C13H28	017312-62-8	50
3	Heptadecane, 2,6,10,15-tetramethyl-			296	C21H44	054833-48-6	43
4	Tetradecane, 4-methyl-			212	C15H32	025117-24-2	43
5	Dodecane, 1-iodo-			296	C12H25I	004292-19-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PPSP-B10-17.5-18.0

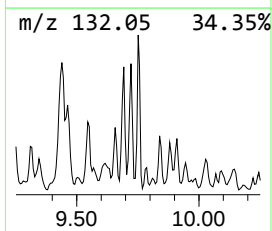
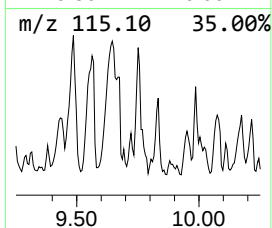
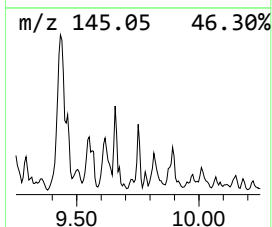
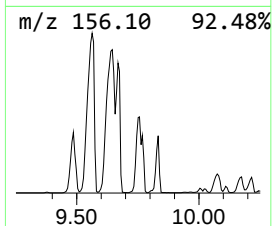
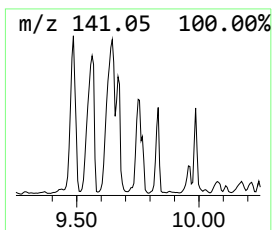
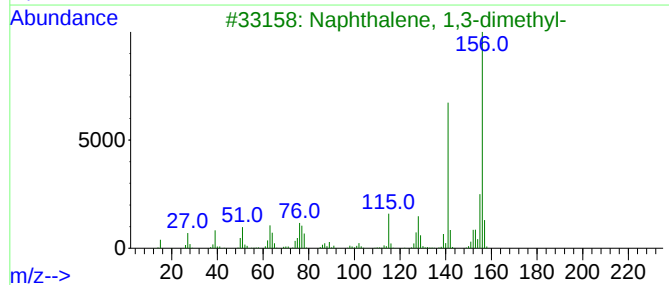
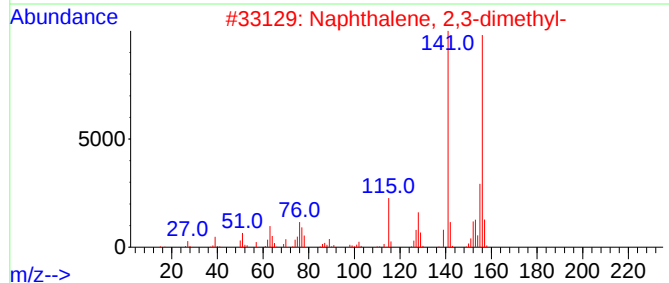
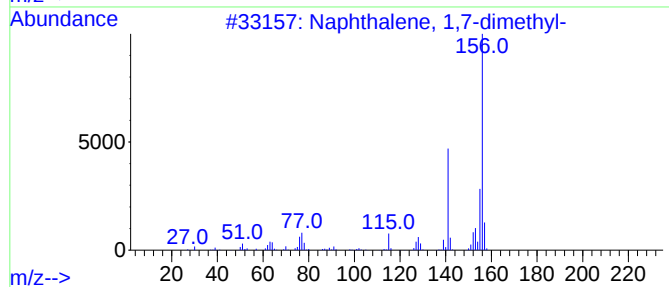
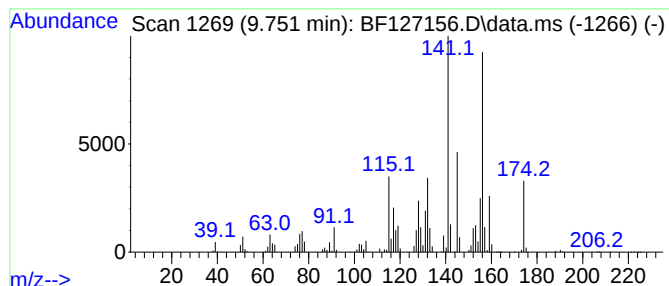
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.751	42.40 ng	11648400	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	91
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	91
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	90
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	90
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

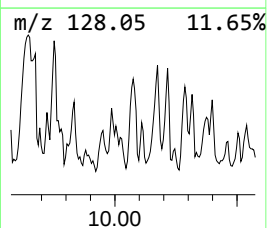
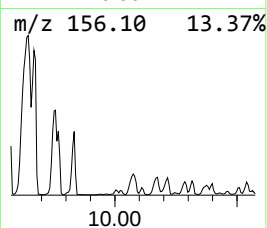
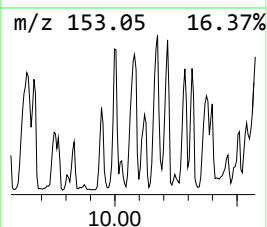
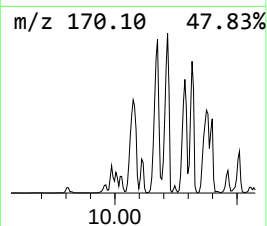
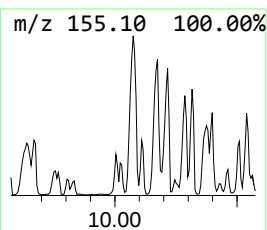
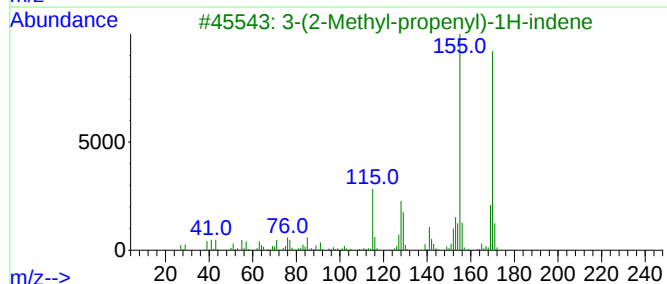
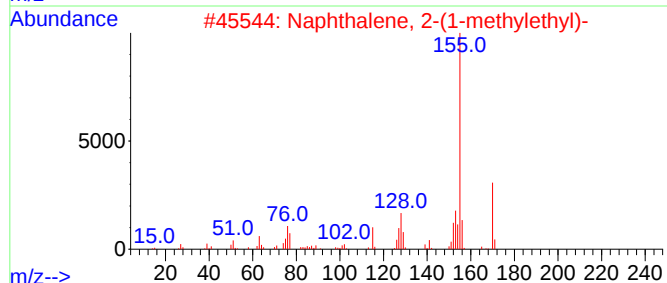
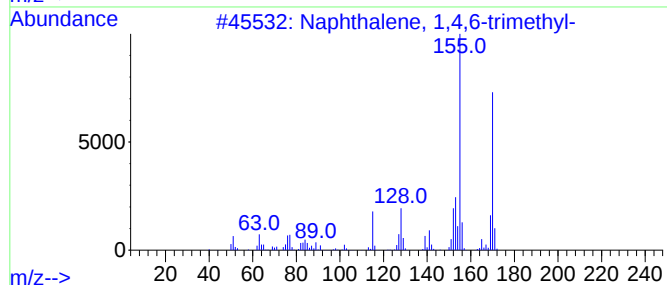
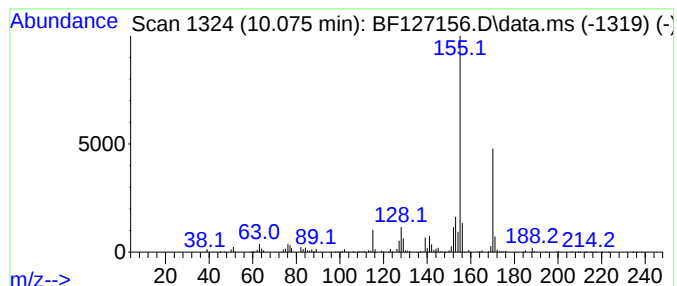
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ClientSampleId :
PPSP-B10-17.5-18.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 1,4,6-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.075	29.78 ng	8180810	Acenaphthene-d10	9.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	91
3	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	90
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	89
5	4a,8a-Methanonaphthalene, 9,9-di...	170	C13H14	038963-97-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PPSP-B10-17.5-18.0

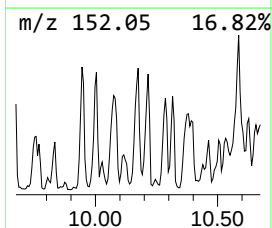
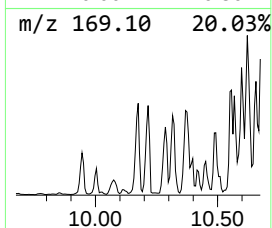
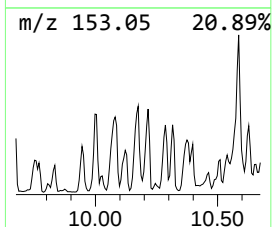
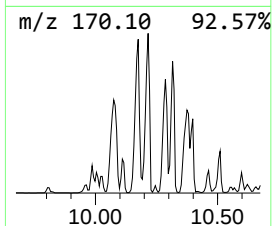
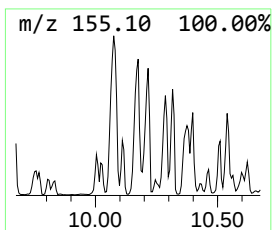
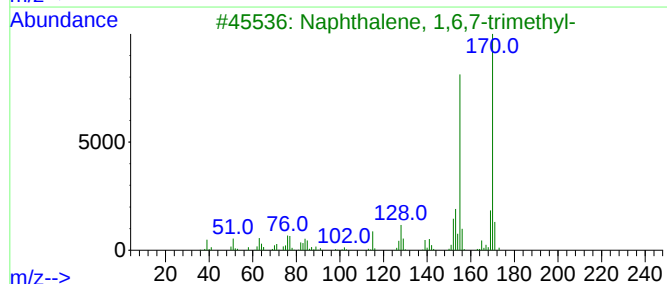
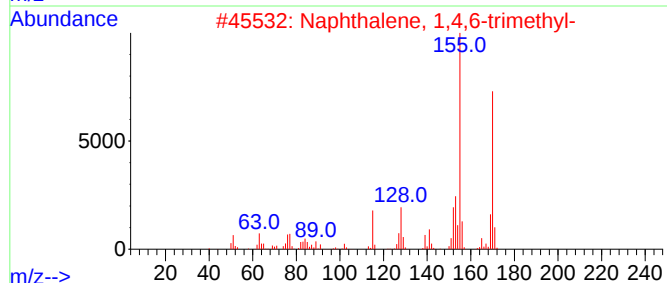
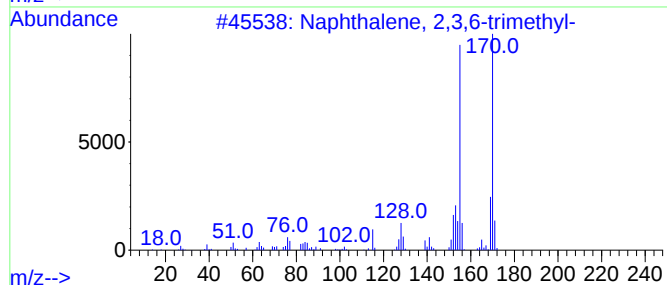
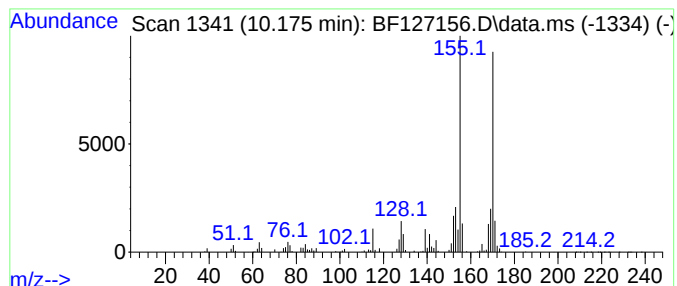
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 2,3,6-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.175	28.76 ng	7902440	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
4		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91
5		1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

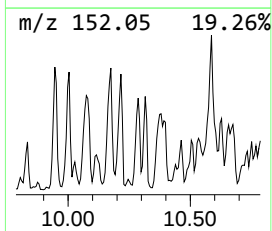
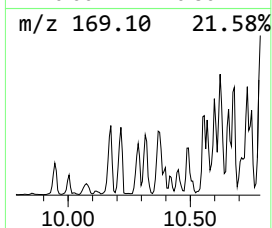
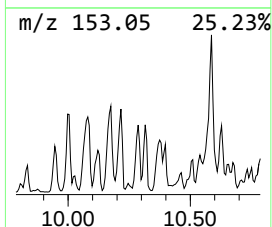
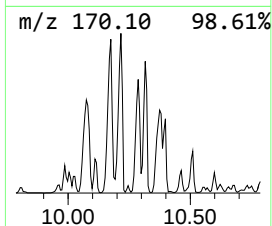
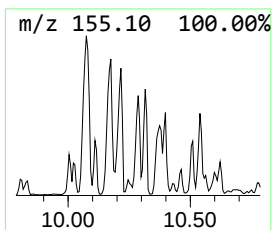
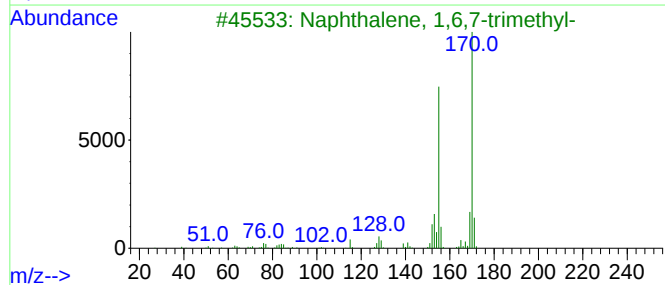
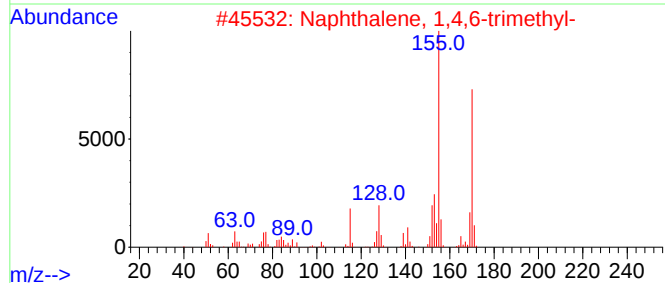
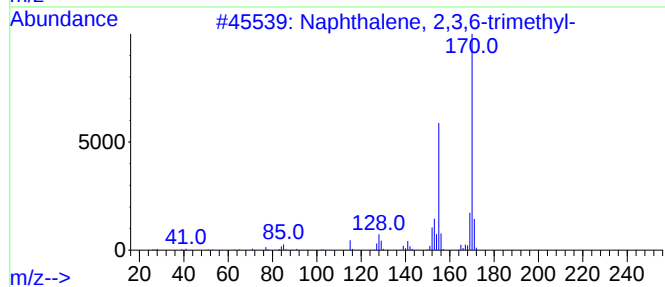
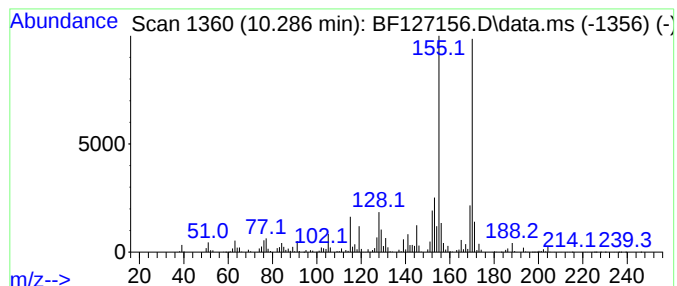
Instrument :
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ClientSampleId :
PPSP-B10-17.5-18.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1,6,7-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.286	27.63 ng	7589970	Acenaphthene-d10	9.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
4	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	91
5	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PPSP-B10-17.5-18.0

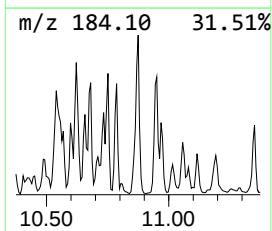
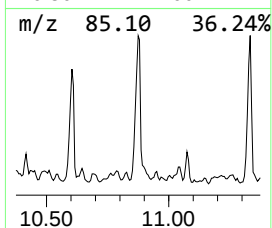
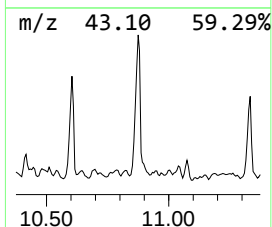
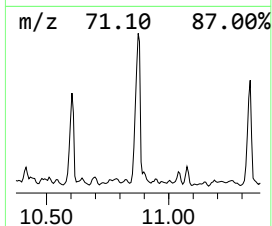
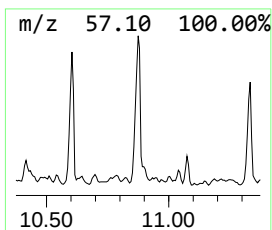
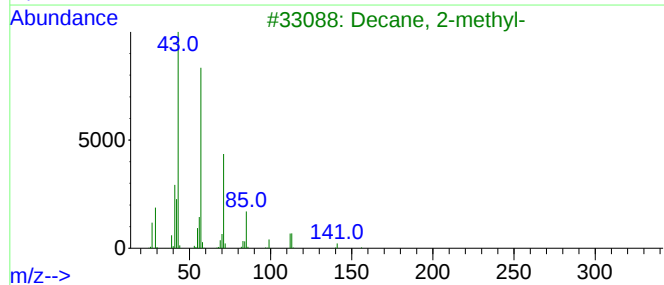
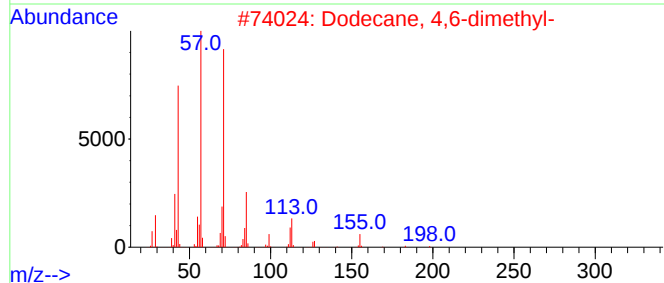
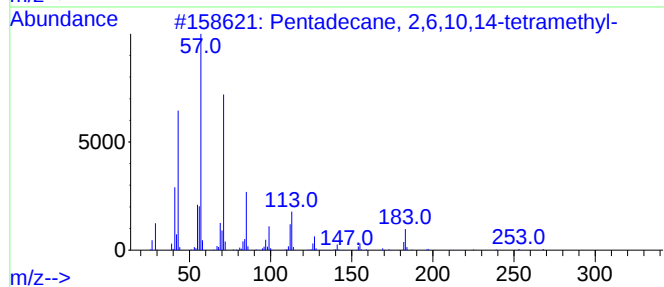
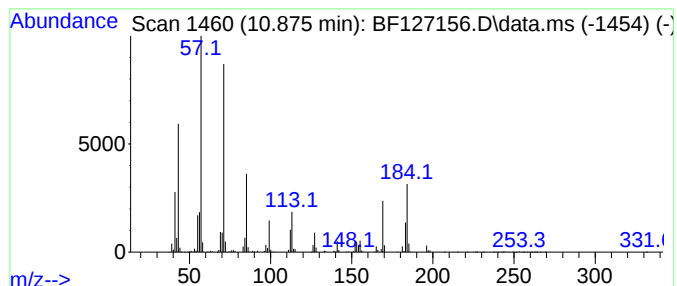
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentadecane, 2,6,10,14-tetr... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.875	90.36 ng	12162300	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72
2			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64
3			Decane, 2-methyl-	156	C11H24	006975-98-0	58
4			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	58
5			Tridecane, 4-methyl-	198	C14H30	026730-12-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PPSP-B10-17.5-18.0

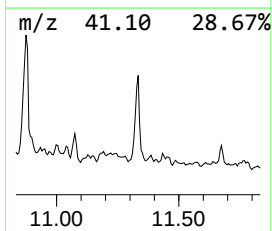
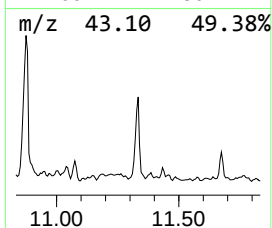
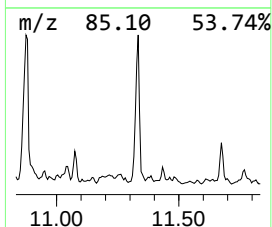
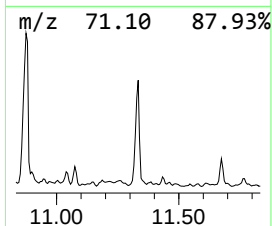
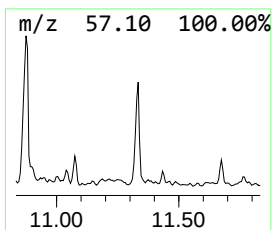
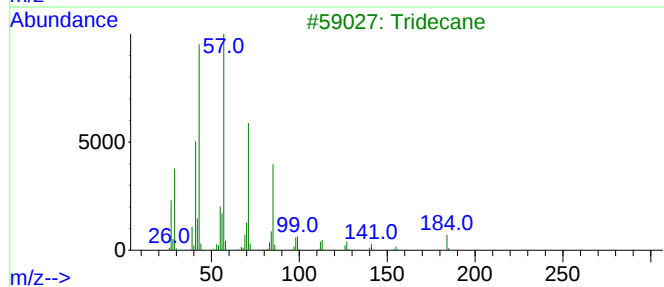
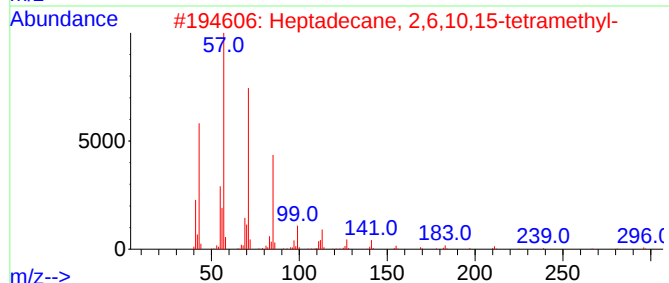
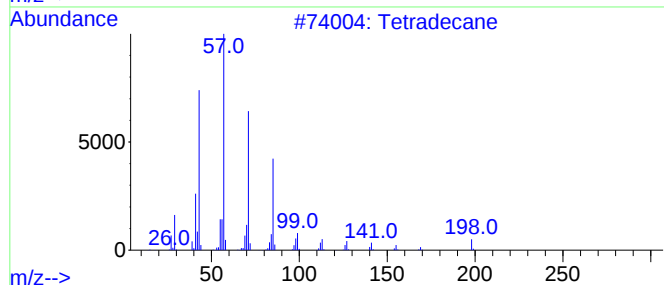
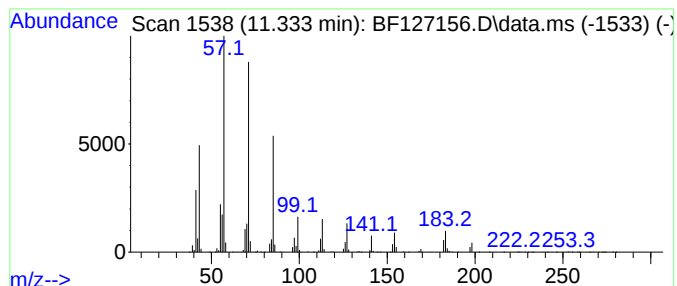
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tetradecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.333	62.76 ng	8447400	Phenanthrene-d10	11.451

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	83
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
3		Tridecane	184	C13H28	000629-50-5	74
4		Hexadecane	226	C16H34	000544-76-3	74
5		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

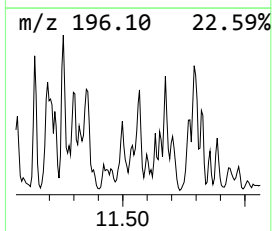
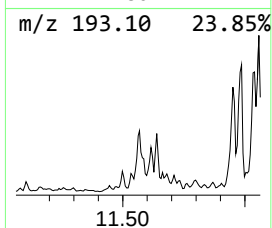
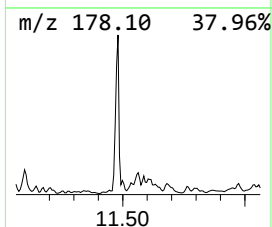
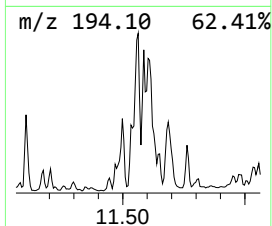
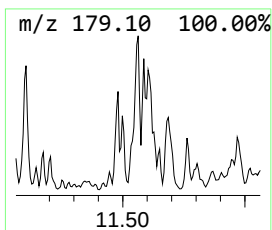
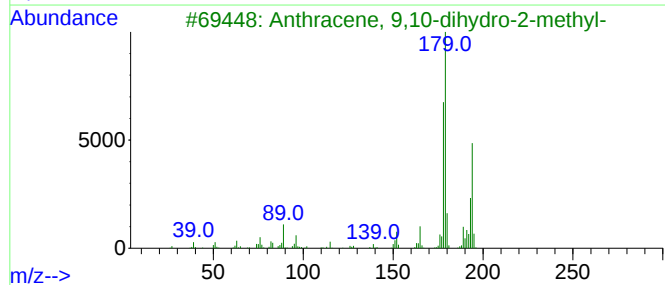
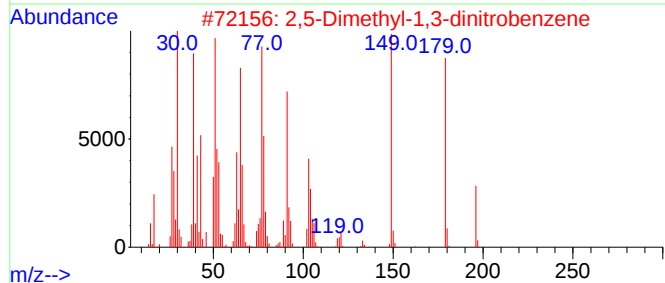
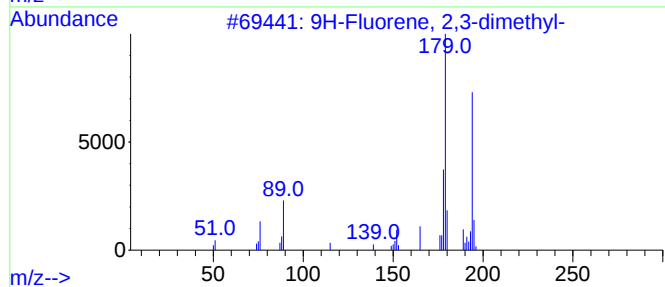
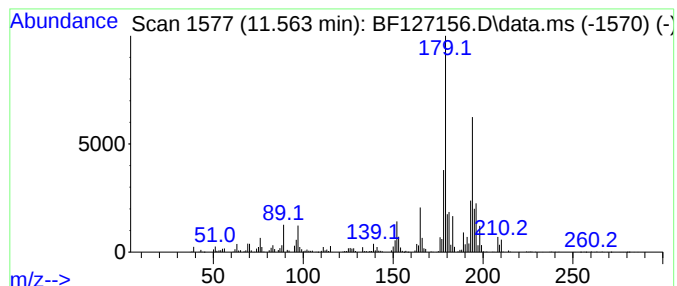
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ClientSampleId :
PPSP-B10-17.5-18.0

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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 9H-Fluorene, 2,3-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.563	25.99 ng	3498410	Phenanthrene-d10		11.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9H-Fluorene, 2,3-dimethyl-		194	C15H14	004612-63-9	93
2	2,5-Dimethyl-1,3-dinitrobenzene		196	C8H8N2O4	1010487-00-2	83
3	Anthracene, 9,10-dihydro-2-methyl-		194	C15H14	000948-67-4	78
4	Benzene, 1-methyl-3-(2-phenyleth...		194	C15H14	014064-48-3	58
5	Phenanthrene, 9,10-dihydro-1-met...		194	C15H14	095676-48-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

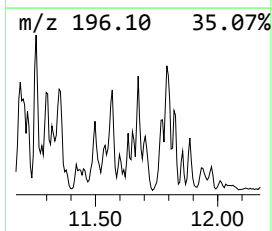
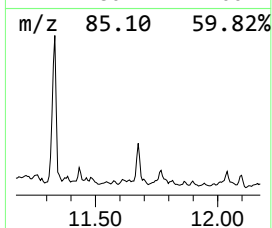
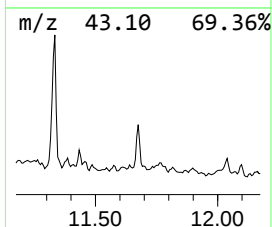
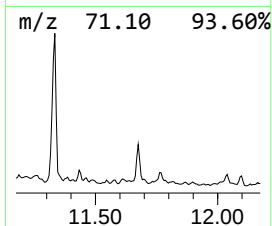
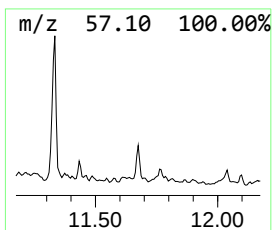
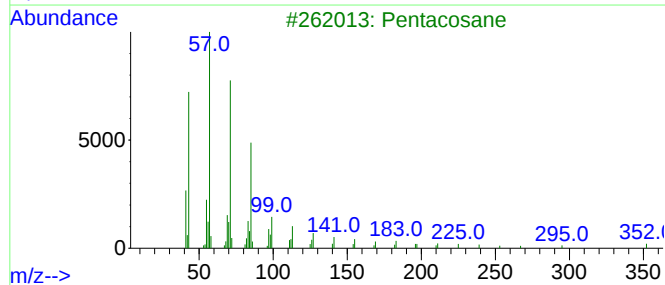
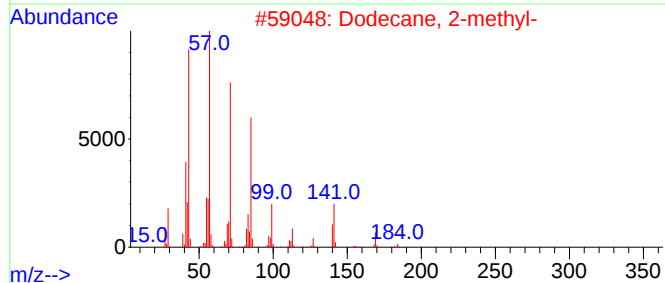
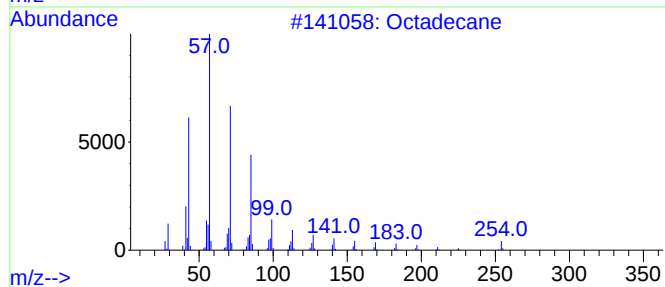
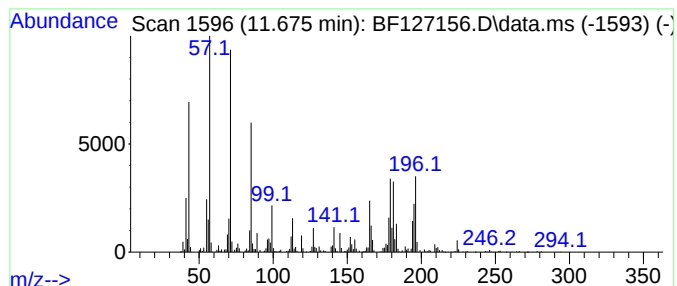
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown11.675 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.675	33.11 ng	4456820	Phenanthrene-d10		11.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	49
2	Dodecane, 2-methyl-		184	C13H28	001560-97-0	49
3	Pentacosane		352	C25H52	000629-99-2	46
4	Tetradecane, 3-methyl-		212	C15H32	018435-22-8	46
5	Heneicosane		296	C21H44	000629-94-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
Data File : BF127156.D
Acq On : 02 Mar 2022 19:35
Operator : CG\JU
Sample : N1708-17
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PPSP-B10-17.5-18.0

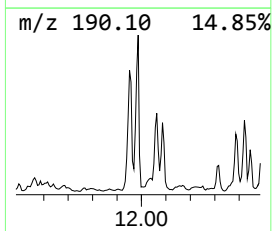
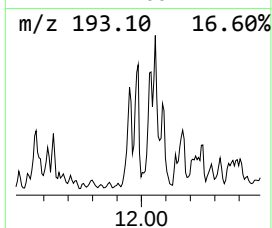
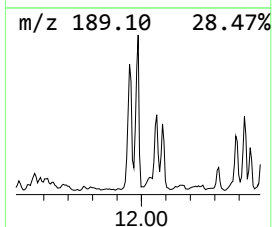
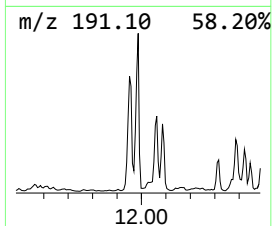
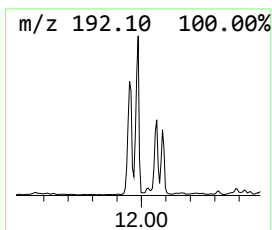
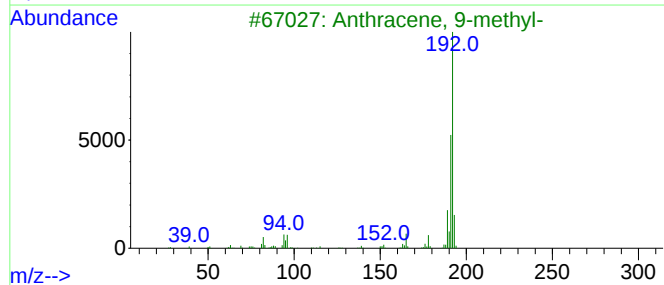
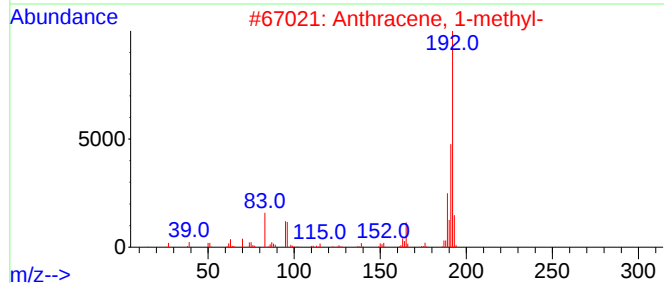
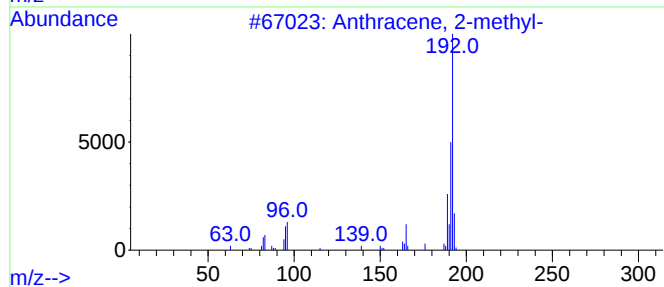
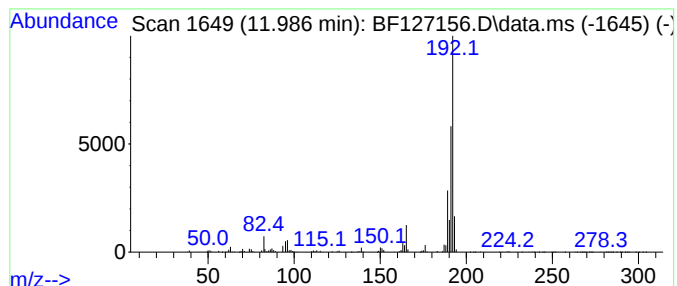
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Anthracene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.986	26.69 ng	3592370	Phenanthrene-d10	11.451

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030222\
 Data File : BF127156.D
 Acq On : 02 Mar 2022 19:35
 Operator : CG\JU
 Sample : N1708-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.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
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p-Mentha-1,5,8-...	7.216	30.6	ng	5070600	1	6.893	3310070	20.0
Benzene, 1-meth...	7.287	32.6	ng	5396200	1	6.893	3310070	20.0
6,6-DIMETHYL-2-...	7.540	29.0	ng	4805110	1	6.893	3310070	20.0
Naphthalene, 1,...	8.710	28.2	ng	9409820	2	8.192	6683890	20.0
Nonane, 3,7-dim...	9.222	23.3	ng	6396910	3	9.963	5494940	20.0
Naphthalene, 1,...	9.439	25.8	ng	7089430	3	9.963	5494940	20.0
Naphthalene, 1-...	9.481	31.9	ng	8776530	3	9.963	5494940	20.0
Naphthalene, 1,...	9.563	51.0	ng	14007000	3	9.963	5494940	20.0
Naphthalene, 2,...	9.639	66.2	ng	18185500	3	9.963	5494940	20.0
Undecane	9.675	44.9	ng	12341700	3	9.963	5494940	20.0
Naphthalene, 1,...	9.751	42.4	ng	11648400	3	9.963	5494940	20.0
Naphthalene, 1,...	10.075	29.8	ng	8180810	3	9.963	5494940	20.0
Naphthalene, 2,...	10.175	28.8	ng	7902440	3	9.963	5494940	20.0
Naphthalene, 1,...	10.286	27.6	ng	7589970	3	9.963	5494940	20.0
Pentadecane, 2,...	10.875	90.4	ng	12162300	4	11.451	2692100	20.0
Tetradecane	11.333	62.8	ng	8447400	4	11.451	2692100	20.0
9H-Fluorene, 2,...	11.563	26.0	ng	3498410	4	11.451	2692100	20.0
unknown11.675	11.675	33.1	ng	4456820	4	11.451	2692100	20.0
Anthracene, 2-m...	11.986	26.7	ng	3592370	4	11.451	2692100	20.0