

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
 Data File : BF114352.D
 Acq On : 17 May 2019 16:20
 Operator : JU/SJ
 Sample : K2871-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04(25.0-27)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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	5.057	502	505	521	rVB2	88745	128469	2.83%	0.147%
2	5.475	572	576	597	rVB	3979724	4035314	88.87%	4.603%
3	6.098	676	682	685	rBV2	155930	199702	4.40%	0.228%
4	6.286	709	714	717	rBV2	79080	94922	2.09%	0.108%
5	6.381	727	730	733	rBV2	191121	195729	4.31%	0.223%
6	6.492	744	749	756	rBV	4418185	4458315	98.18%	5.085%
7	6.586	761	765	767	rBV2	70944	104178	2.29%	0.119%
8	6.622	767	771	774	rVV	4638191	4162529	91.67%	4.748%
9	6.845	805	809	816	rVB2	1532077	1492929	32.88%	1.703%
10	7.004	831	836	839	rVV	3524605	3237116	71.29%	3.692%
11	7.122	853	856	859	rVB	252197	252502	5.56%	0.288%
12	7.151	859	861	864	rBV	117363	108488	2.39%	0.124%
13	7.186	864	867	870	rVV2	109764	119680	2.64%	0.137%
14	7.233	870	875	879	rVV3	196742	273905	6.03%	0.312%
15	7.269	879	881	886	rVB5	73381	98741	2.17%	0.113%
16	7.381	893	900	901	rBV3	202547	320868	7.07%	0.366%
17	7.410	901	905	914	rVB	2716176	2998654	66.04%	3.420%
18	7.486	914	918	922	rVB5	192472	304604	6.71%	0.347%
19	7.551	922	929	932	rBV3	153090	349039	7.69%	0.398%
20	7.598	934	937	939	rVV2	109790	124580	2.74%	0.142%
21	7.622	939	941	944	rVV2	199098	209155	4.61%	0.239%
22	7.651	944	946	949	rVB	225660	164269	3.62%	0.187%
23	7.745	954	962	964	rBV4	253181	347537	7.65%	0.396%
24	7.769	964	966	969	rVB3	325549	273812	6.03%	0.312%
25	7.804	969	972	975	rBV3	222414	277350	6.11%	0.316%
26	7.839	975	978	981	rBV2	164875	147205	3.24%	0.168%
27	7.875	981	984	989	rBV6	252007	336591	7.41%	0.384%
28	7.922	989	992	994	rVB	169888	122459	2.70%	0.140%
29	7.945	994	996	1002	rVB5	133819	160607	3.54%	0.183%
30	7.998	1002	1005	1008	rBV3	141198	204354	4.50%	0.233%
31	8.033	1008	1011	1014	rBV	104008	88446	1.95%	0.101%
32	8.063	1014	1016	1017	rBV2	116200	78898	1.74%	0.090%
33	8.080	1017	1019	1021	rVV3	68313	83027	1.83%	0.095%
34	8.128	1021	1027	1033	rVV	1799945	2036079	44.84%	2.322%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	8.186	1033	1037	1040	rVB	997715	946503	20.84%	1.080%
36	8.216	1040	1042	1045	rBV3	260017	233110	5.13%	0.266%
37	8.280	1050	1053	1055	rBV4	152489	173968	3.83%	0.198%
38	8.304	1055	1057	1059	rVV2	107862	101239	2.23%	0.115%
39	8.333	1059	1062	1064	rVB	352910	282587	6.22%	0.322%
40	8.416	1069	1076	1083	rBV5	646096	1159323	25.53%	1.322%
41	8.469	1083	1085	1088	rVB2	253901	228598	5.03%	0.261%
42	8.510	1088	1092	1095	rVB4	285728	330570	7.28%	0.377%
43	8.551	1095	1099	1101	rBV	1164133	1121335	24.69%	1.279%
44	8.569	1101	1102	1104	rVB	235720	134025	2.95%	0.153%
45	8.592	1104	1106	1109	rVB3	238906	193118	4.25%	0.220%
46	8.633	1110	1113	1115	rBV	352307	278179	6.13%	0.317%
47	8.669	1115	1119	1121	rBV4	433089	458609	10.10%	0.523%
48	8.716	1125	1127	1131	rBV4	301631	355579	7.83%	0.406%
49	8.757	1131	1134	1136	rBV3	150239	163673	3.60%	0.187%
50	8.816	1141	1144	1147	rVB4	627609	626976	13.81%	0.715%
51	8.875	1151	1154	1156	rBV4	308993	266747	5.87%	0.304%
52	8.910	1156	1160	1162	rBV4	248666	402133	8.86%	0.459%
53	9.004	1173	1176	1178	rBV4	312151	388107	8.55%	0.443%
54	9.033	1178	1181	1185	rVB4	570551	665806	14.66%	0.759%
55	9.075	1185	1188	1192	rBV4	169322	266765	5.87%	0.304%
56	9.122	1193	1196	1198	rVV2	310373	234646	5.17%	0.268%
57	9.151	1198	1201	1205	rBV	1546791	1466367	32.29%	1.673%
58	9.198	1206	1209	1214	rVB	3962611	3877185	85.38%	4.422%
59	9.239	1214	1216	1218	rBV2	125351	97504	2.15%	0.111%
60	9.286	1220	1224	1228	rBV2	762716	1117366	24.61%	1.274%
61	9.398	1241	1243	1247	rVB	409117	404588	8.91%	0.461%
62	9.469	1251	1255	1258	rVV2	524989	692015	15.24%	0.789%
63	9.539	1264	1267	1269	rBV	851836	702271	15.47%	0.801%
64	9.604	1273	1278	1281	rBV2	2259102	2623885	57.78%	2.993%
65	9.627	1281	1282	1284	rVB	136863	79882	1.76%	0.091%
66	9.657	1285	1287	1289	rBV	389483	285224	6.28%	0.325%
67	9.751	1299	1303	1306	rBV3	498418	634385	13.97%	0.724%
68	9.792	1307	1310	1312	rBV	625547	524300	11.55%	0.598%
69	9.880	1318	1325	1328	rBV2	1918105	2561036	56.40%	2.921%
70	9.916	1329	1331	1334	rVB3	409212	338418	7.45%	0.386%
71	9.986	1340	1343	1347	rBV	673256	793791	17.48%	0.905%

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72	10.039	1348	1352	1354	rBV5	282454	462750	10.19%	0.528%
73	10.074	1355	1358	1363	rBV2	710774	1009253	22.23%	1.151%
74	10.139	1365	1369	1376	rVB7	788425	1440210	31.72%	1.643%
75	10.198	1376	1379	1382	rBV2	816125	934045	20.57%	1.065%
76	10.227	1382	1384	1386	rBV	437833	337109	7.42%	0.385%
77	10.251	1386	1388	1391	rVB2	400192	382463	8.42%	0.436%
78	10.292	1391	1395	1396	rBV3	812405	845261	18.61%	0.964%
79	10.427	1415	1418	1424	rBV5	556359	974957	21.47%	1.112%
80	10.533	1431	1436	1440	rVB	2245918	2466435	54.32%	2.813%
81	10.580	1442	1444	1449	rVB4	449805	592067	13.04%	0.675%
82	10.627	1449	1452	1453	rBV2	265859	292022	6.43%	0.333%
83	10.651	1453	1456	1457	rBV	398138	295145	6.50%	0.337%
84	10.674	1457	1460	1463	rBV	2258437	2000911	44.06%	2.282%
85	10.798	1476	1481	1484	rBV	4159745	4540855	100.00%	5.179%
86	11.010	1515	1517	1520	rVB2	679669	548062	12.07%	0.625%
87	11.263	1557	1560	1566	rVB	3139666	3518874	77.49%	4.014%
88	11.374	1575	1579	1581	rBV	1948083	2004506	44.14%	2.286%
89	11.398	1581	1583	1586	rVB	824322	718744	15.83%	0.820%
90	11.616	1617	1620	1623	rBV2	1186904	1139039	25.08%	1.299%
91	11.874	1662	1664	1666	rBV	629854	596515	13.14%	0.680%
92	11.986	1680	1683	1685	rVB2	955193	852823	18.78%	0.973%
93	12.351	1742	1745	1753	rVB3	901585	1180003	25.99%	1.346%
94	12.427	1756	1758	1761	rVB	888527	741758	16.34%	0.846%
95	12.592	1783	1786	1789	rBV	1022366	948301	20.88%	1.082%
96	12.821	1822	1825	1826	rBV	686057	589394	12.98%	0.672%
97	12.874	1832	1834	1838	rVB	600743	636175	14.01%	0.726%
98	12.957	1845	1848	1851	rVB	2654920	2434215	53.61%	2.776%
99	14.033	2028	2031	2033	rVB2	1022447	1047568	23.07%	1.195%
100	15.539	2285	2287	2294	rVB	1046321	1345141	29.62%	1.534%

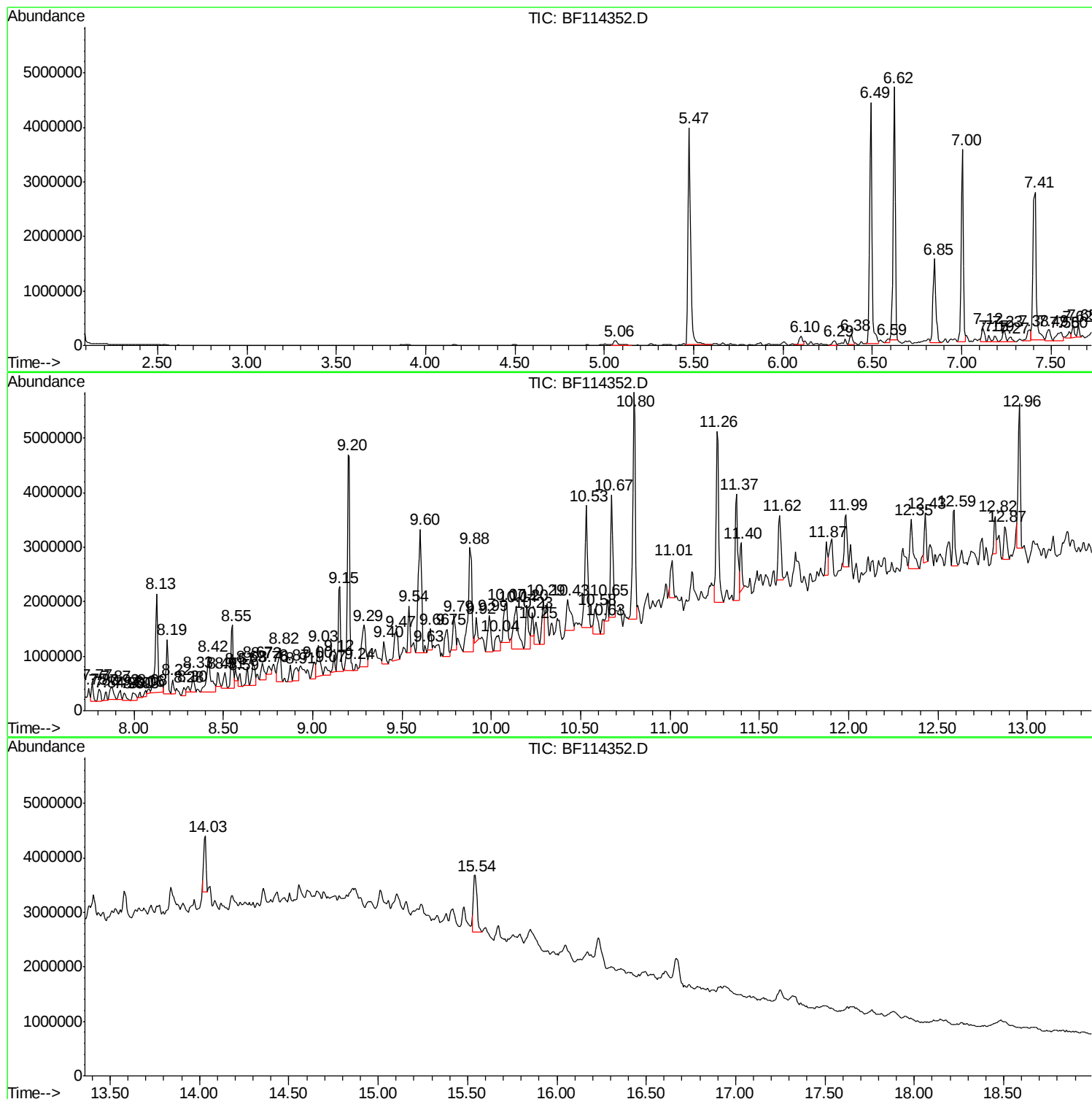
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Instrument :
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ClientSampled :
SB-04(25.0-27)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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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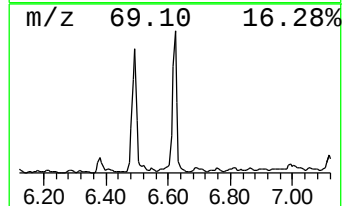
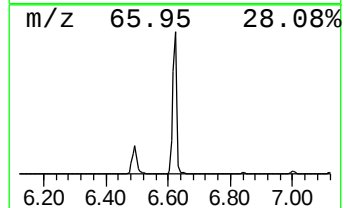
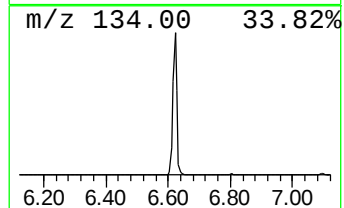
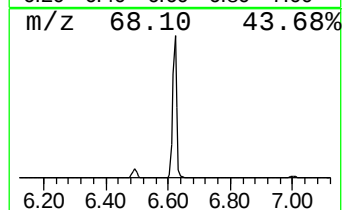
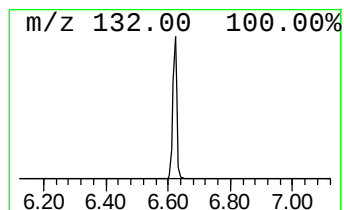
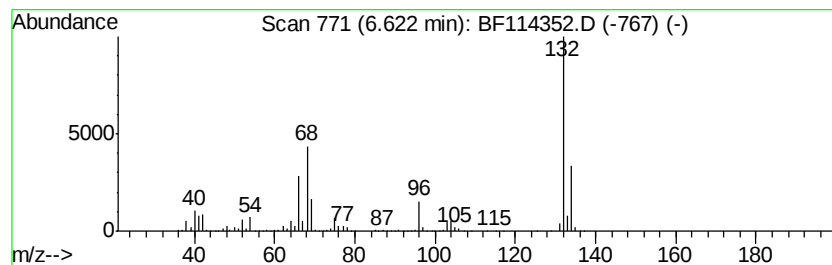
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	55.76 ng	4162530	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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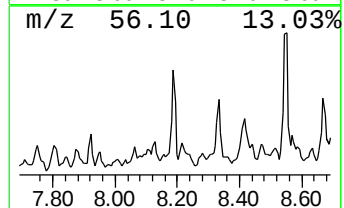
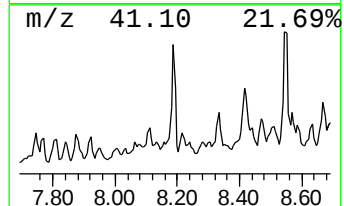
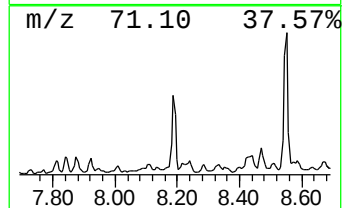
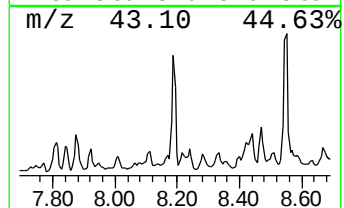
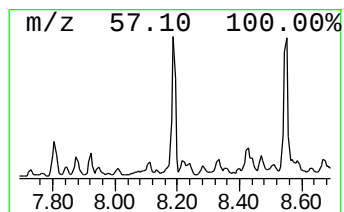
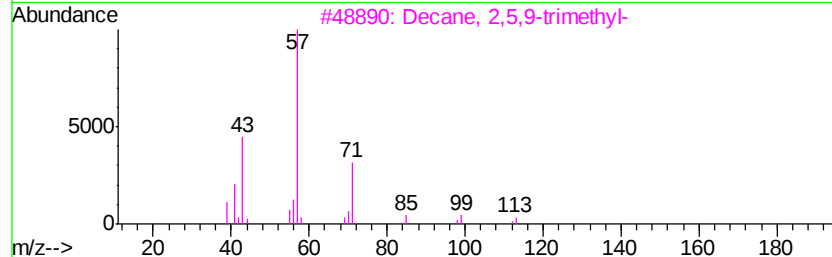
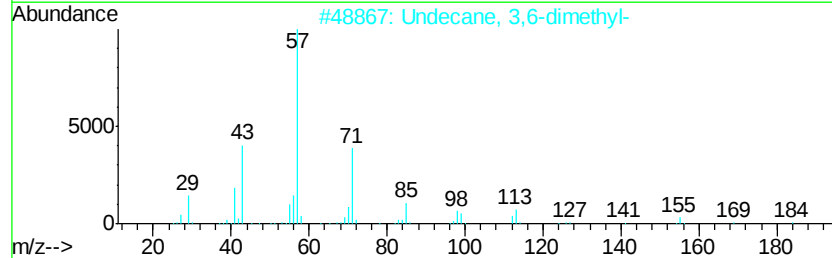
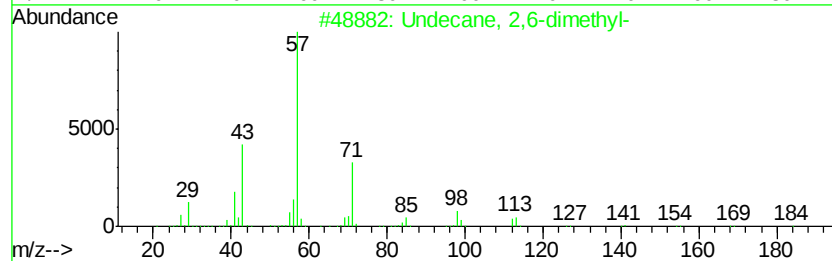
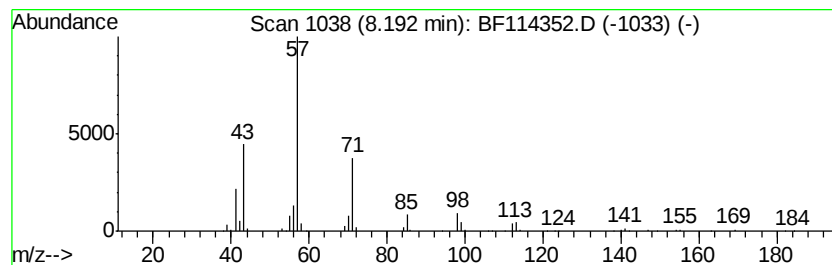
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Peak Number 3 Undecane, 2,6-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.19	9.30 ng	946503	Naphthalene-d8	8.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	97
2	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	78
3	Decane, 2,5,9-trimethyl-	184	C13H28	062108-22-9	72
4	Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	59
5	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	59



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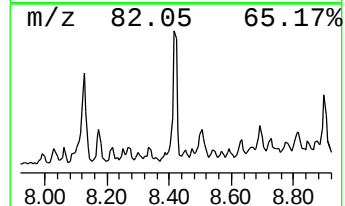
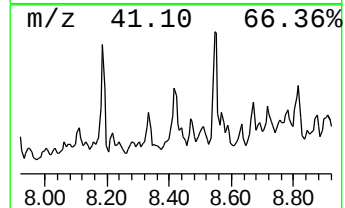
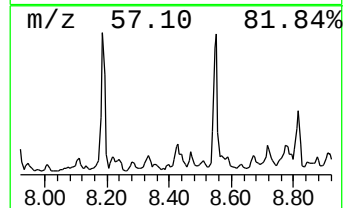
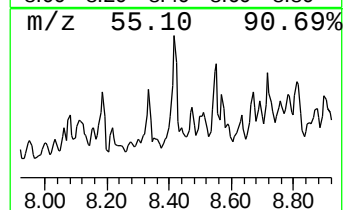
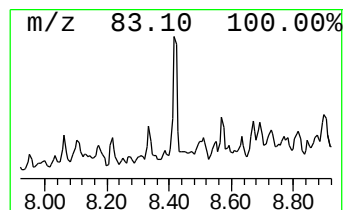
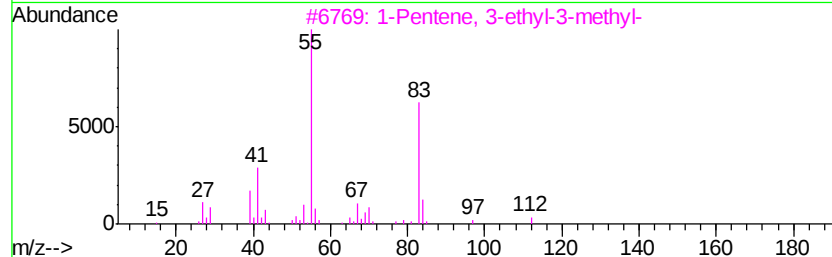
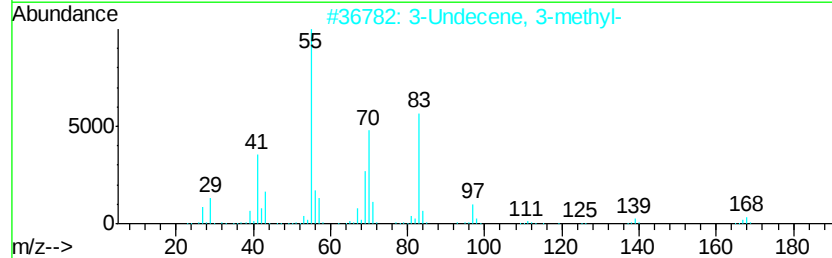
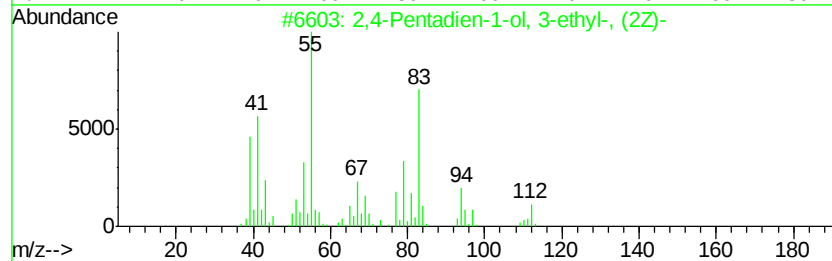
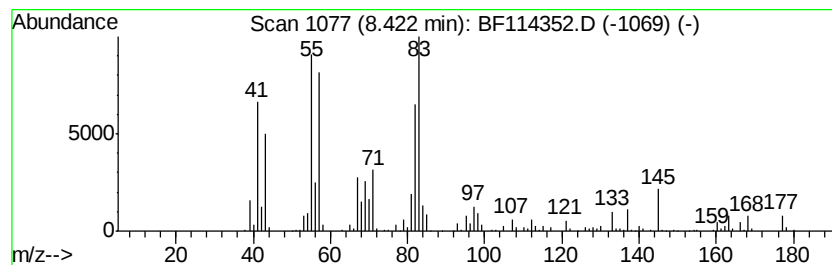
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown8.42 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	11.39 ng	1159320	Napthalene-d8	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Pentadien-1-ol, 3-ethyl-, (2Z)-	112	C7H12O	1000142-17-6	46		
2	3-Undecene, 3-methyl-	168	C12H24	023381-94-4	43		
3	1-Pentene, 3-ethyl-3-methyl-	112	C8H16	006196-60-7	38		
4	3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	38		
5	2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	38		



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
Sample : K2871-02
Misc :
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Instrument :
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ClientSampleId :
SB-04(25.0-27)

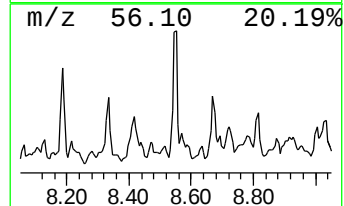
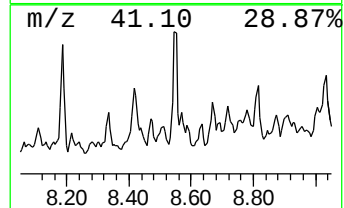
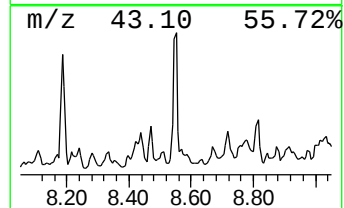
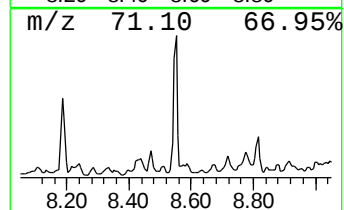
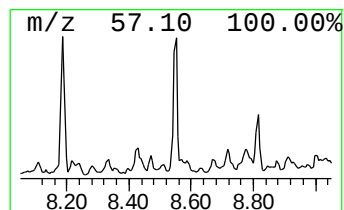
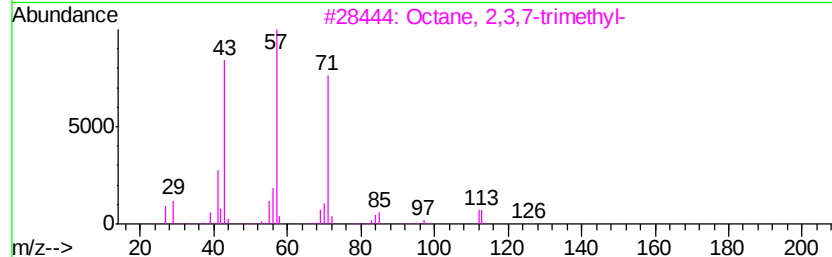
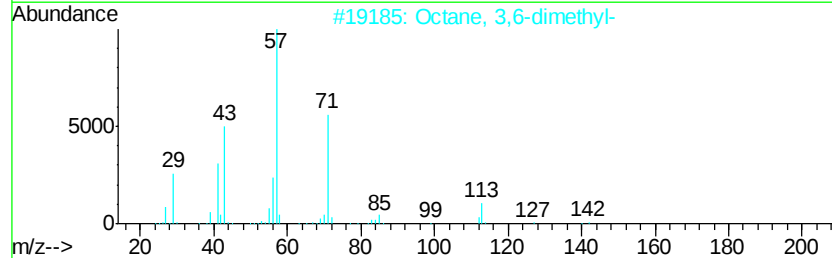
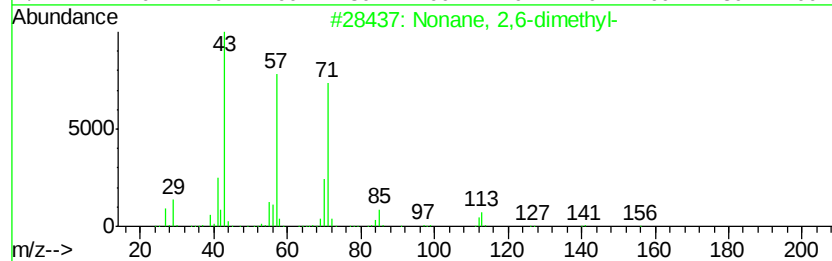
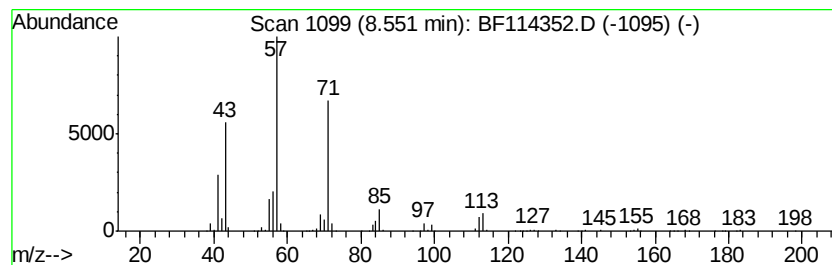
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonane, 2,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.55	11.01 ng	1121340	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	80
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
3		Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
4		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
5		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
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ClientSampleId :
SB-04(25.0-27)

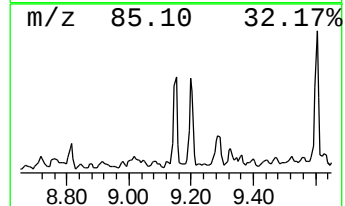
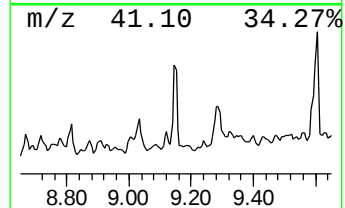
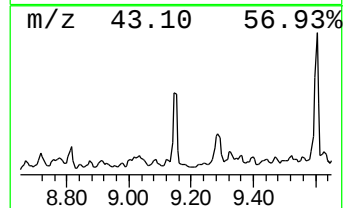
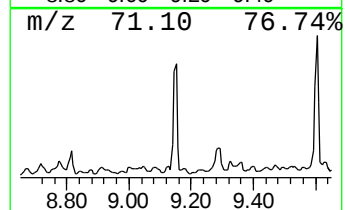
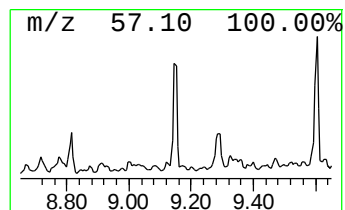
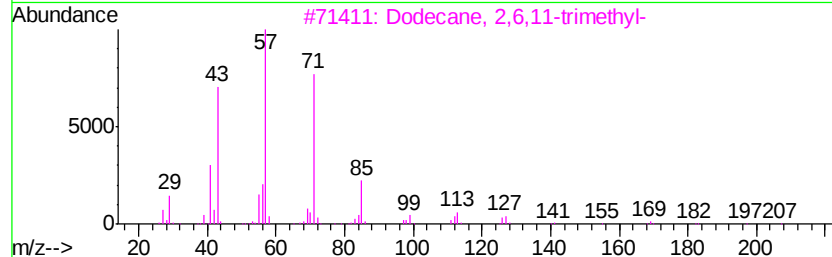
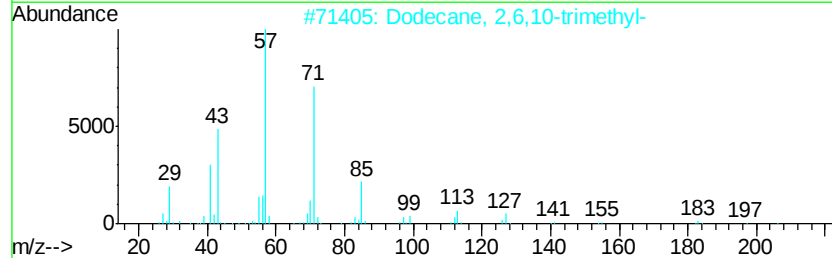
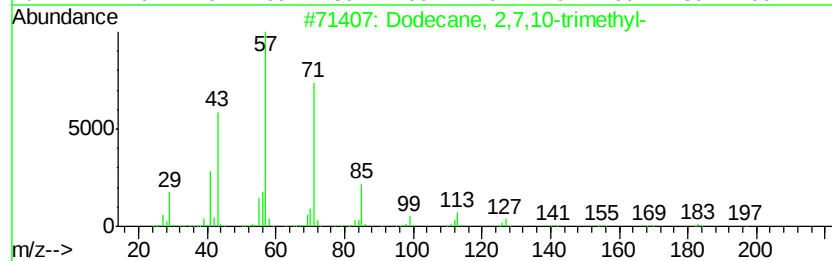
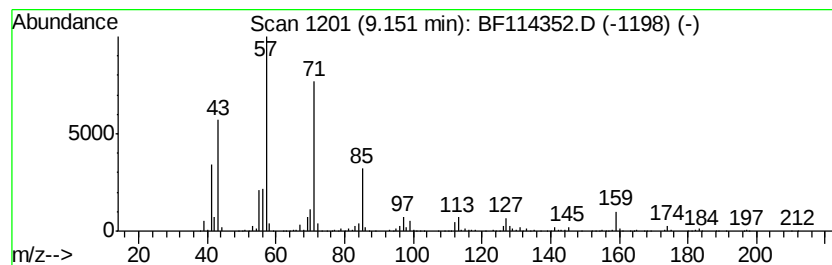
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dodecane, 2,7,10-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.15	11.45 ng	1466370	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	87
2		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	80
3		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80
4		Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
5		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
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ClientSampleId :
SB-04(25.0-27)

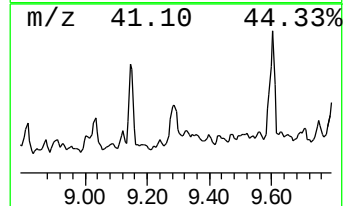
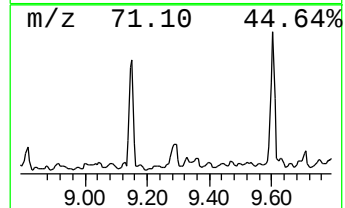
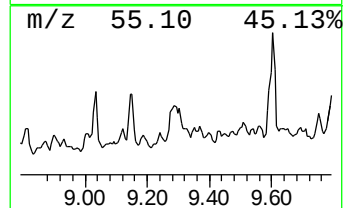
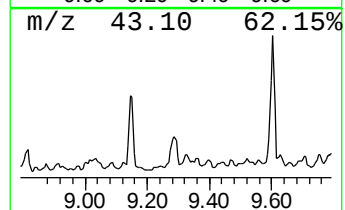
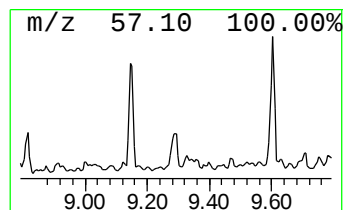
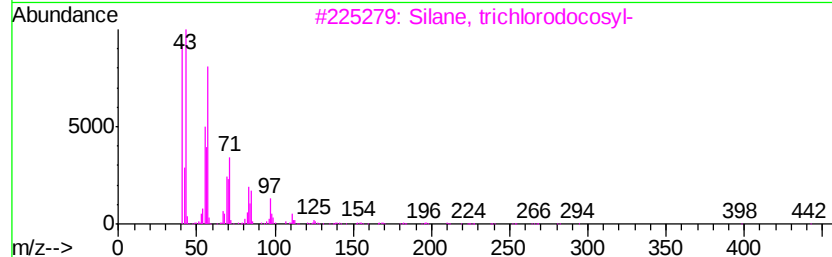
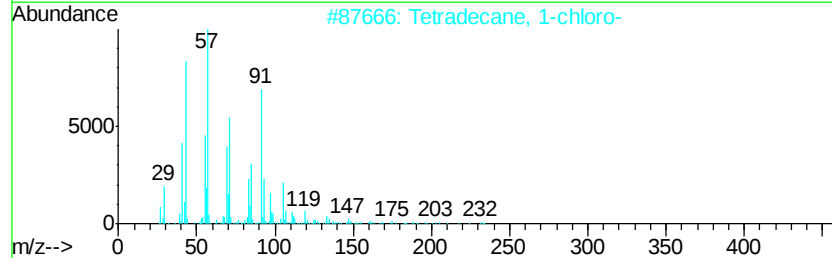
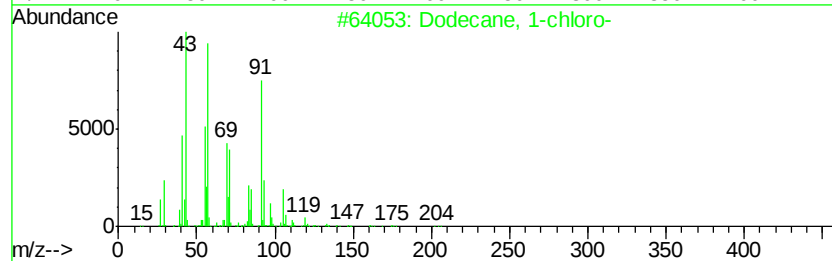
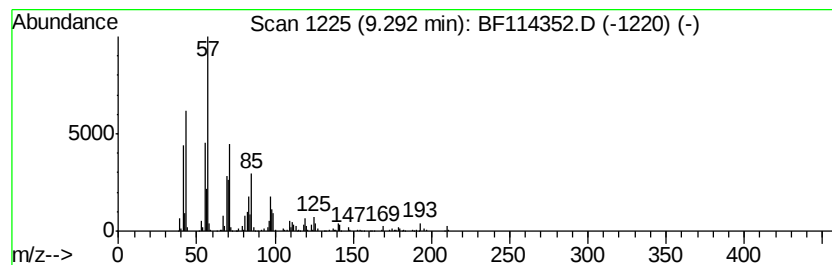
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 1-chloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.29	8.73 ng	1117370	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 1-chloro-	204	C12H25Cl	000112-52-7	74
2		Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	64
3		Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	58
4		Hexadecane, 1-bromo-	304	C16H33Br	000112-82-3	58
5		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
Sample : K2871-02
Misc :
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Instrument :
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ClientSampleId :
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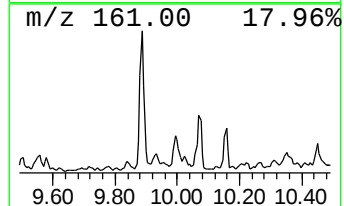
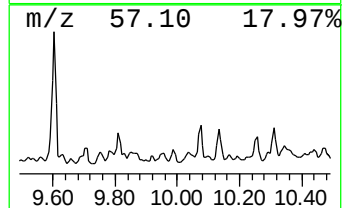
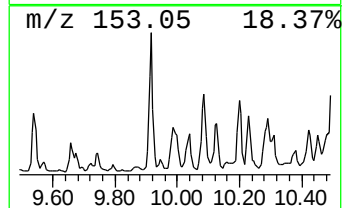
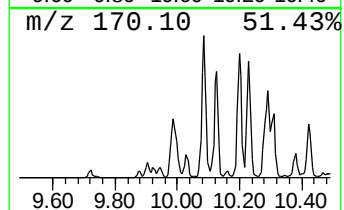
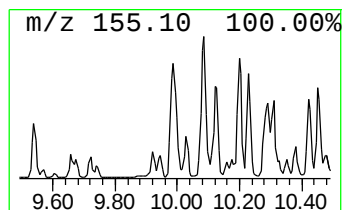
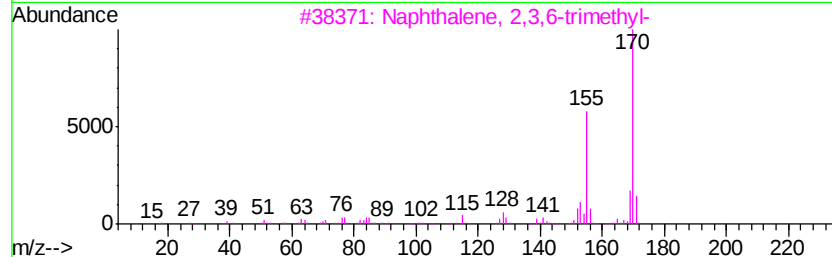
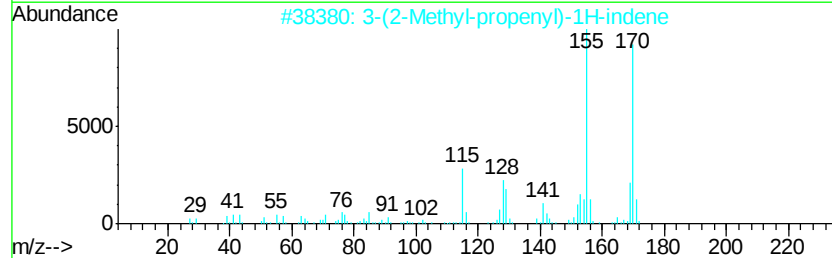
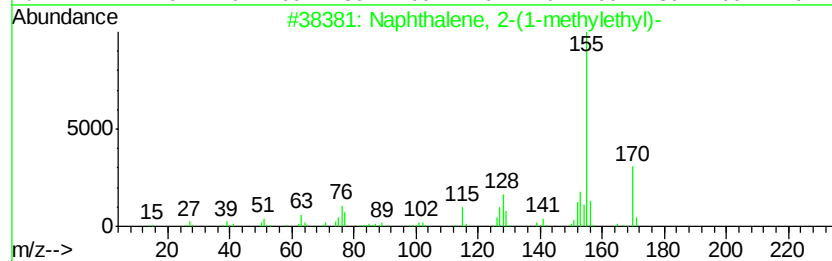
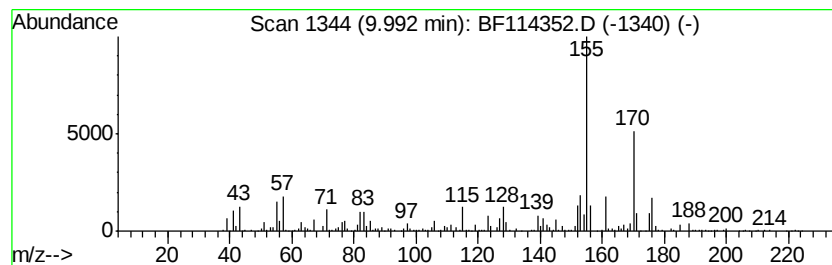
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-(1-methyleth... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.99	6.20 ng	793791	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	95
2		3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	74
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	70
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	70
5		Ethanone, 1-(5-chloro-2-hydroxyp...	170	C8H7ClO2	001450-74-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
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Operator : JU/SJ
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Instrument :
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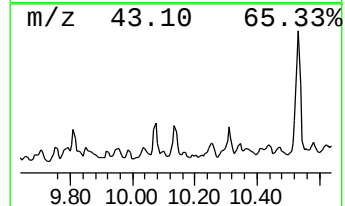
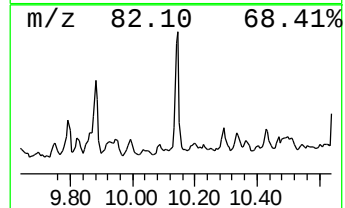
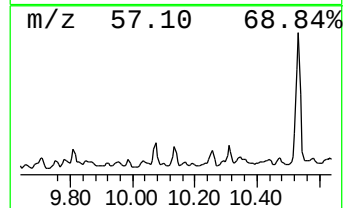
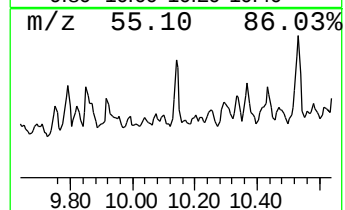
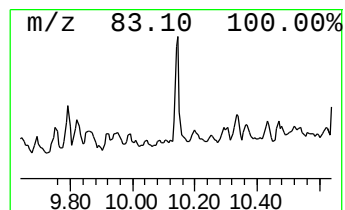
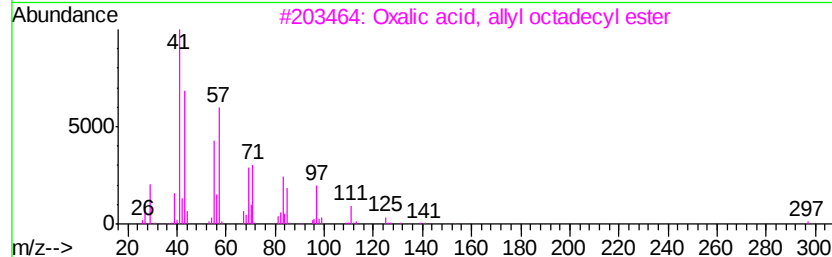
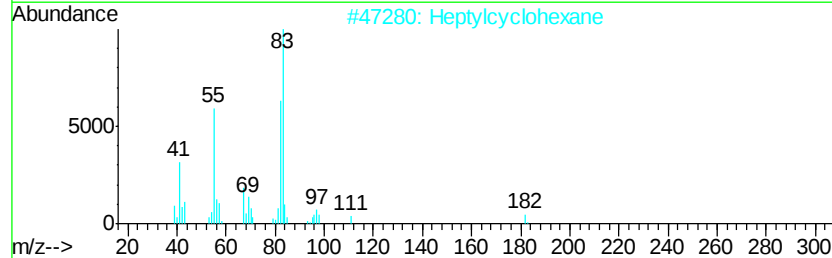
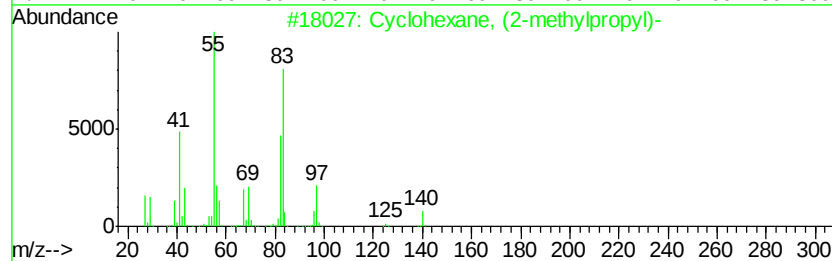
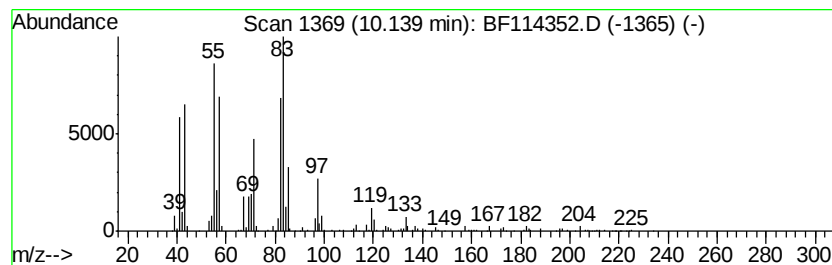
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclohexane, (2-methylpropyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	11.25 ng	1440210	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	62
2		Heptylcyclohexane	182	C13H26	005617-41-4	49
3		Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	47
4		Propanoic acid, 2,2-dimethyl-, c...	184	C11H20O2	029878-49-7	43
5		Oxalic acid, dicyclohexyl ester	254	C14H22O4	1000309-30-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
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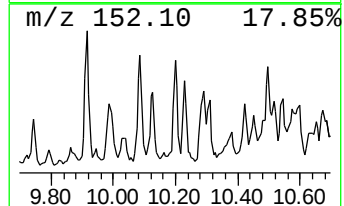
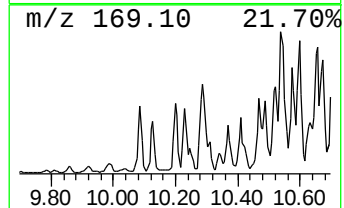
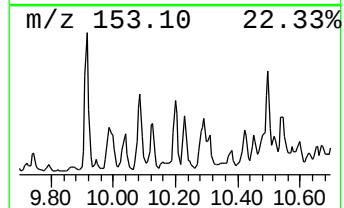
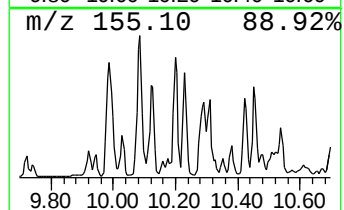
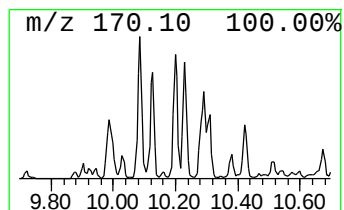
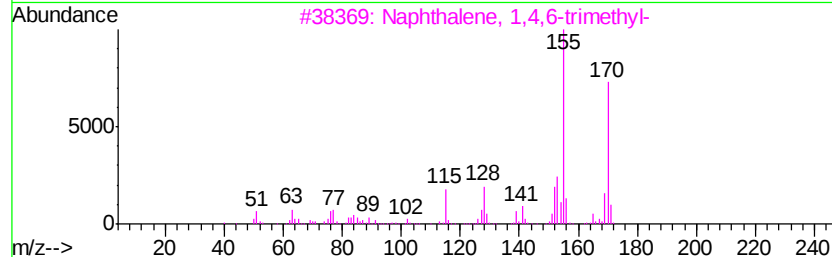
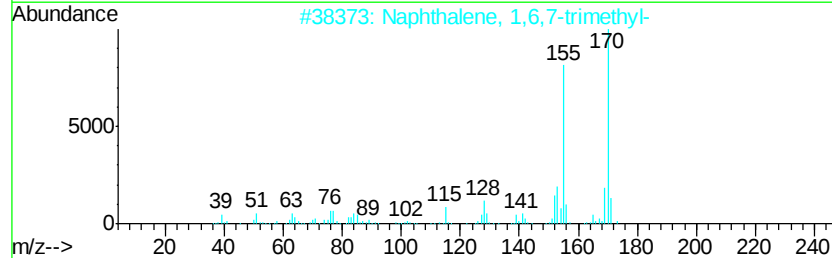
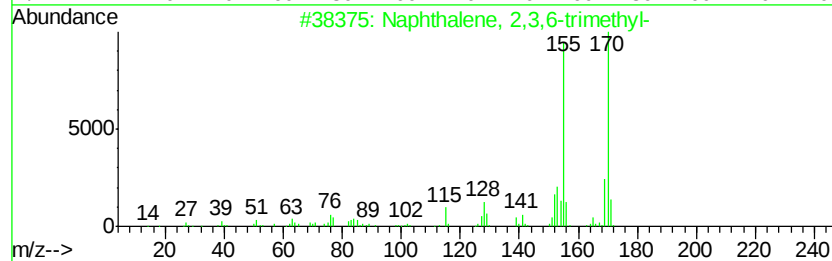
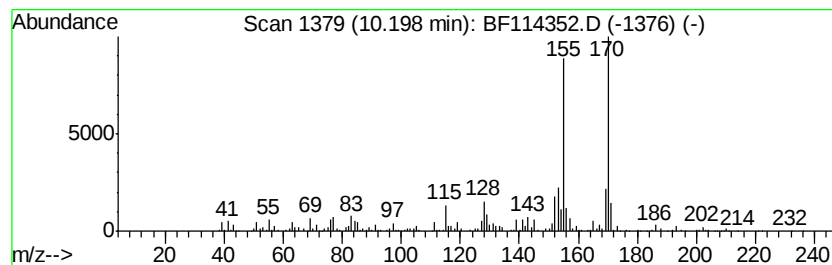
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2,3,6-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.20	7.29 ng	934045	Acenaphthene-d10	9.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	98
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	94
5			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
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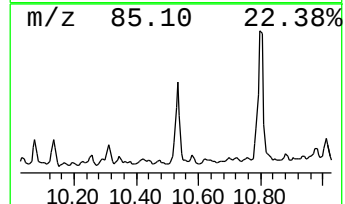
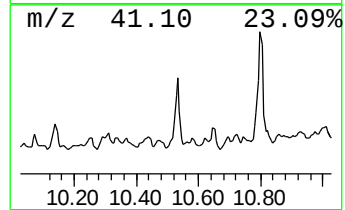
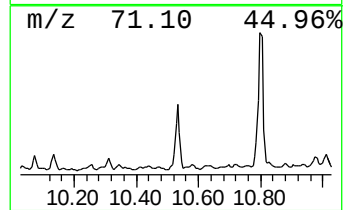
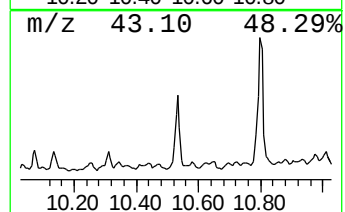
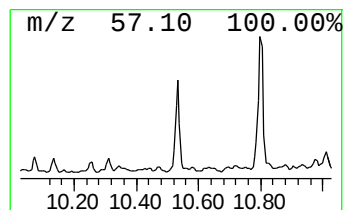
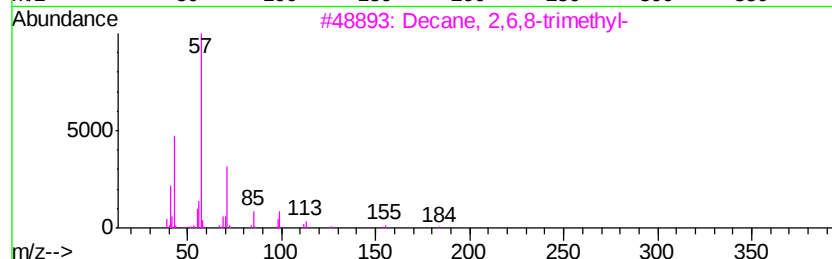
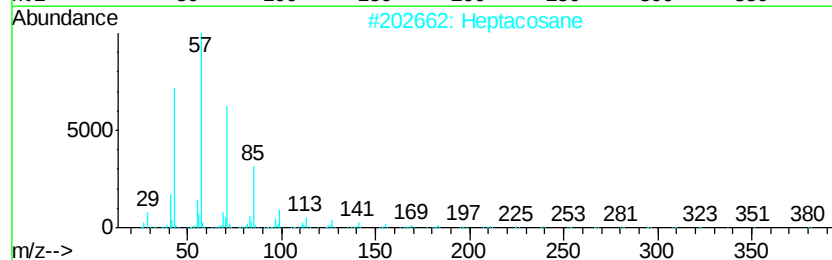
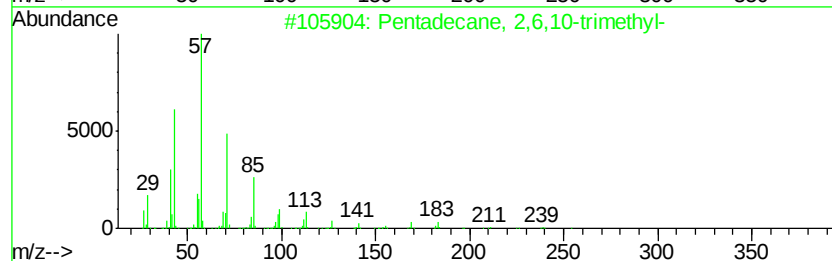
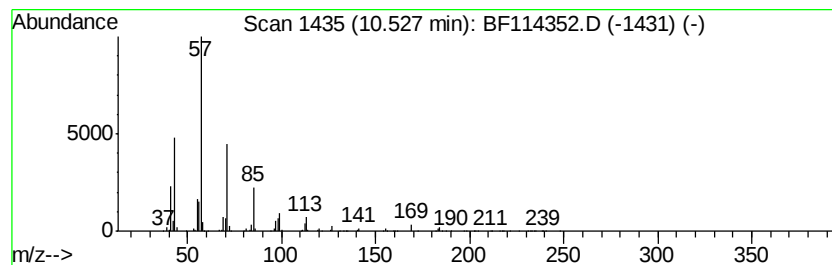
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 Pentadecane, 2,6,10-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	19.26 ng	2466440	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Heptacosane	380	C27H56	000593-49-7	86
3		Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	76
4		Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	76
5		Heneicosane	296	C21H44	000629-94-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
Sample : K2871-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-04(25.0-27)

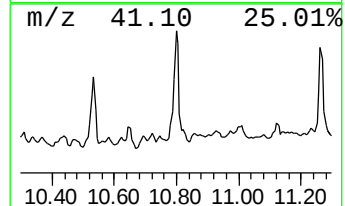
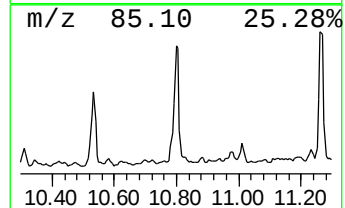
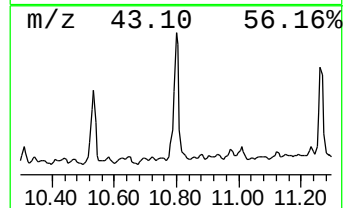
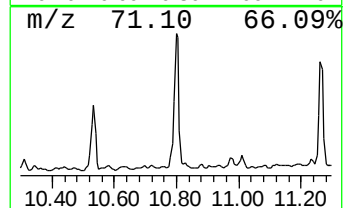
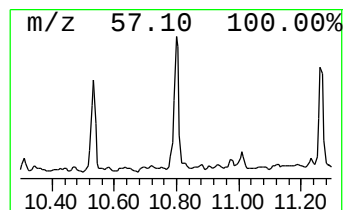
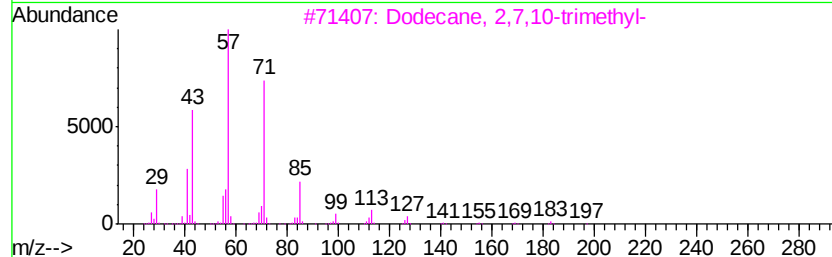
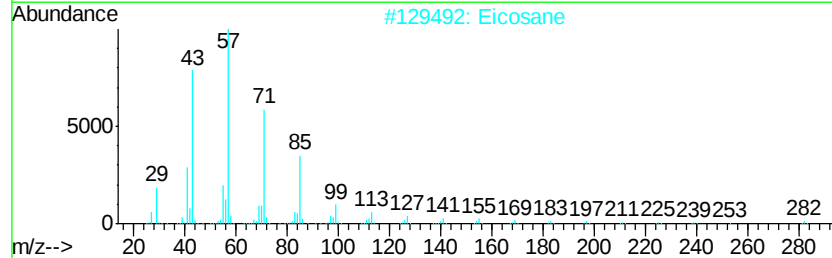
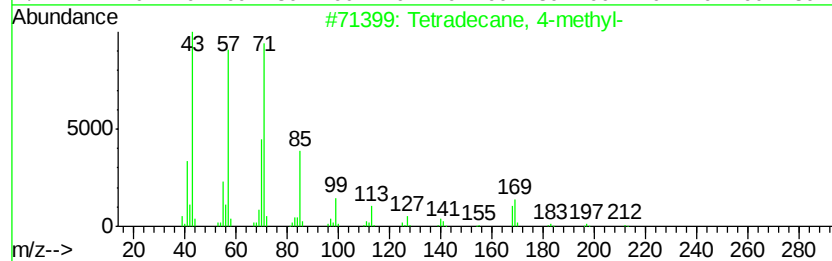
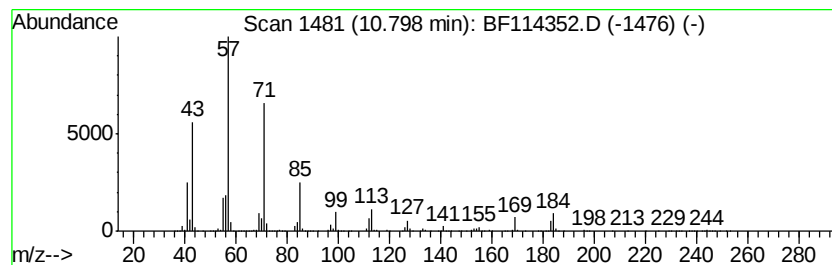
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetradecane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	45.31 ng	4540860	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 4-methyl-	212	C15H32	025117-24-2	72
2		Eicosane	282	C20H42	000112-95-8	72
3		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	58
4		Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	58
5		Tridecane, 3-methyl-	198	C14H30	006418-41-3	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
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Operator : JU/SJ
Sample : K2871-02
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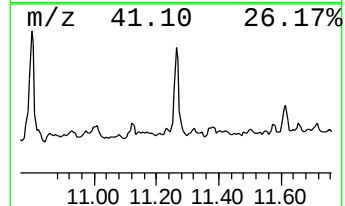
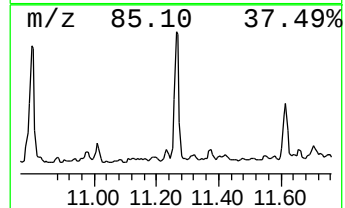
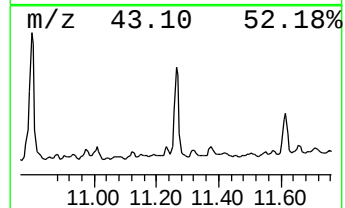
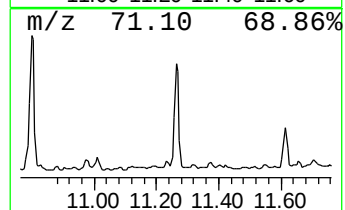
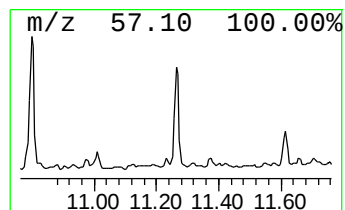
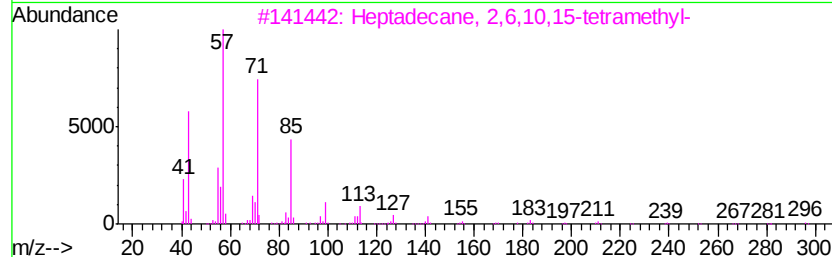
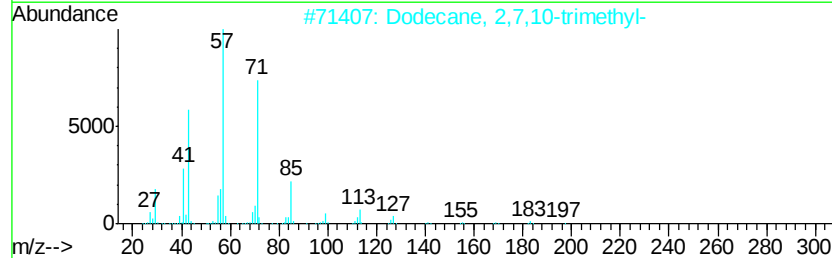
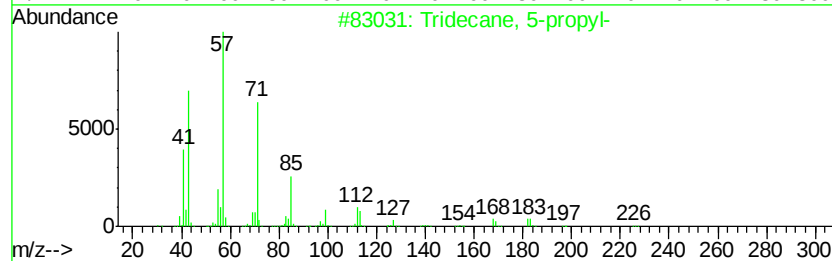
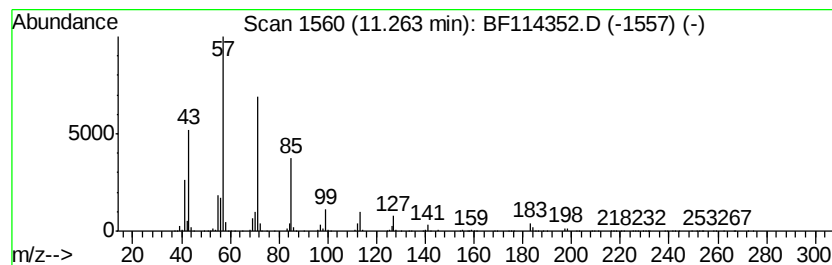
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tridecane, 5-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	35.11 ng	3518870	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	87
2		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	78
5		Tritetracontane	605	C43H88	007098-21-7	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
Sample : K2871-02
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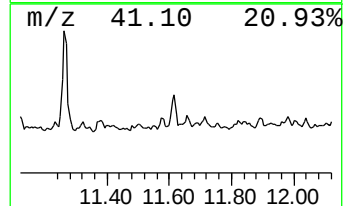
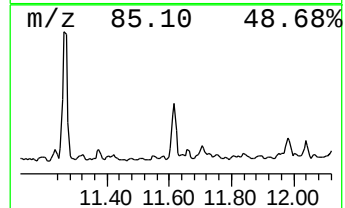
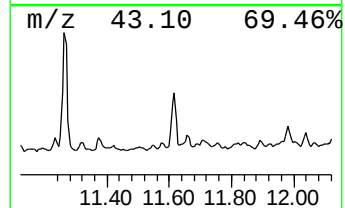
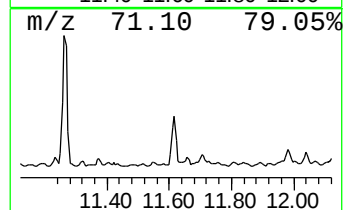
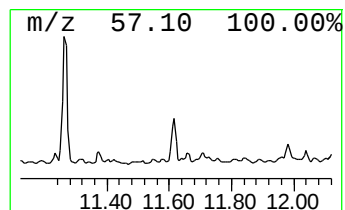
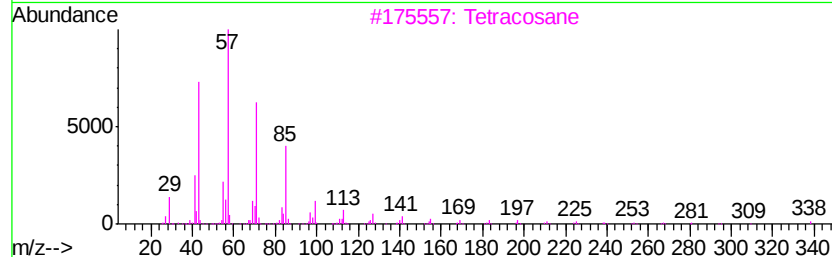
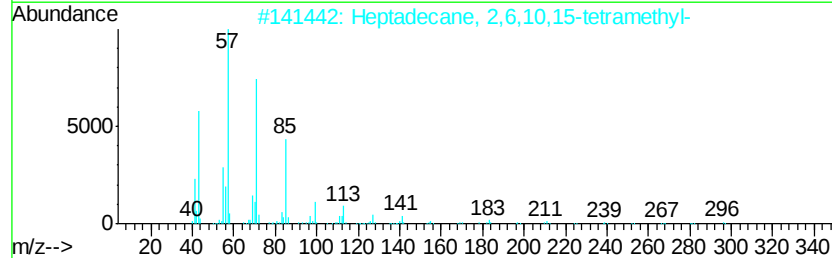
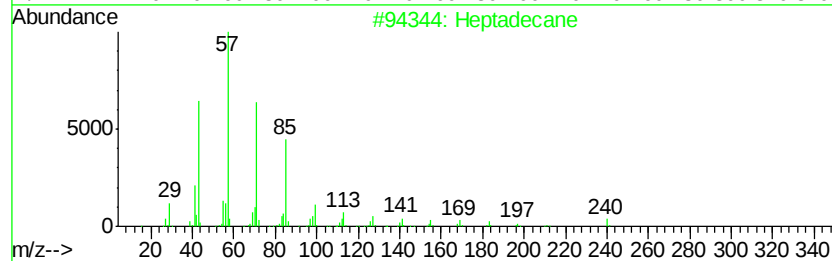
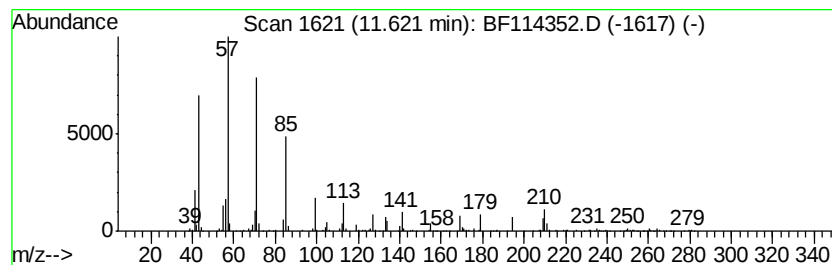
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	11.36 ng	1139040	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	90
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
3		Tetracosane	338	C24H50	000646-31-1	86
4		Heptadecane, 4-methyl-	254	C18H38	026429-11-8	83
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
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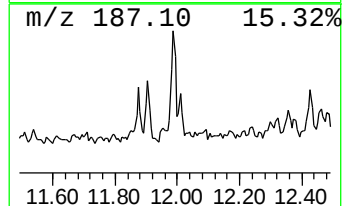
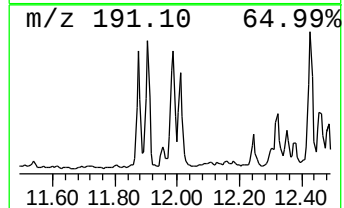
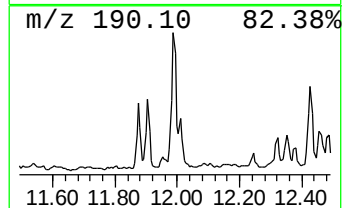
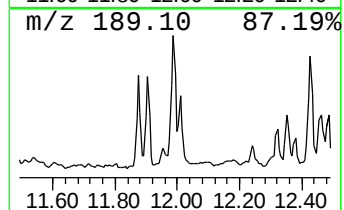
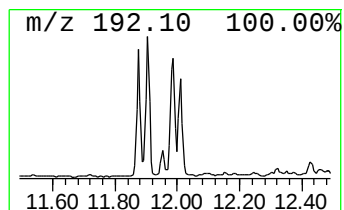
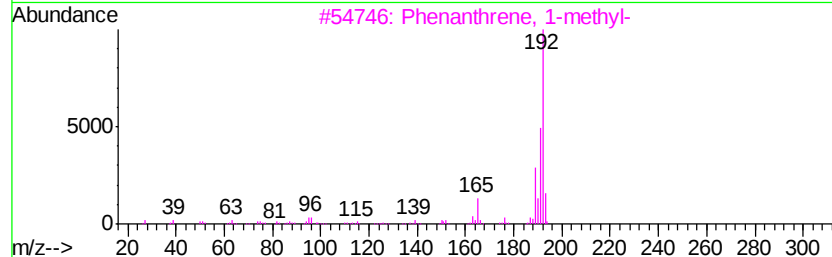
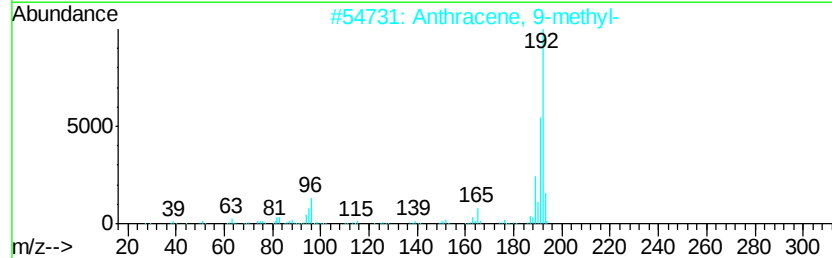
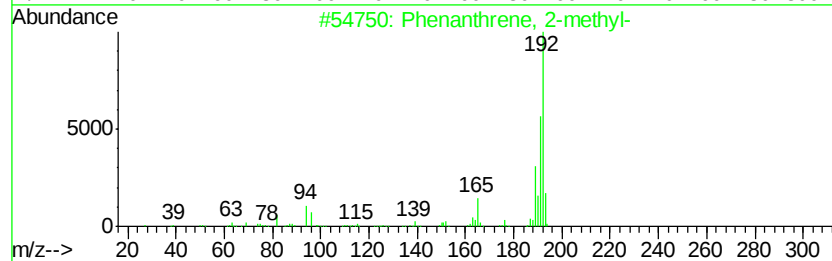
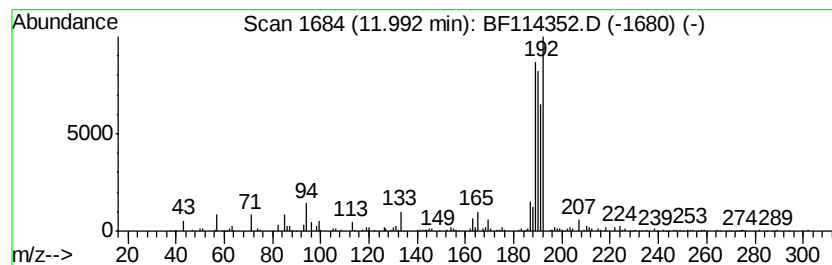
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Phenanthrene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	8.51 ng	852823	Phenanthrene-d10	11.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	64
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	64
4			1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	64
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
Operator : JU/SJ
Sample : K2871-02
Misc :
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ClientSampleId :
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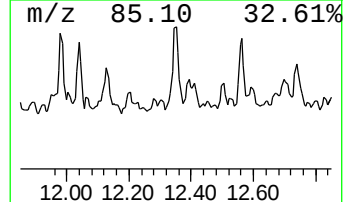
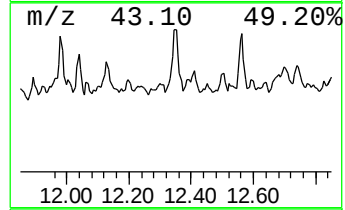
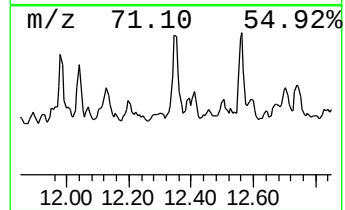
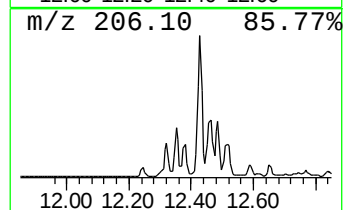
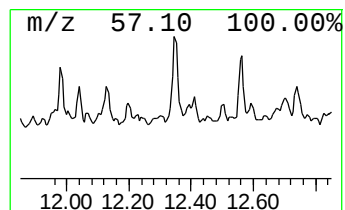
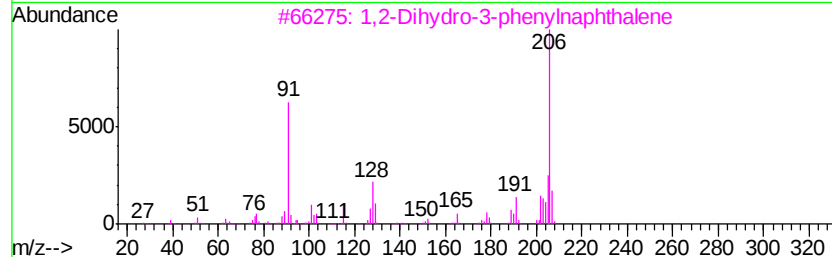
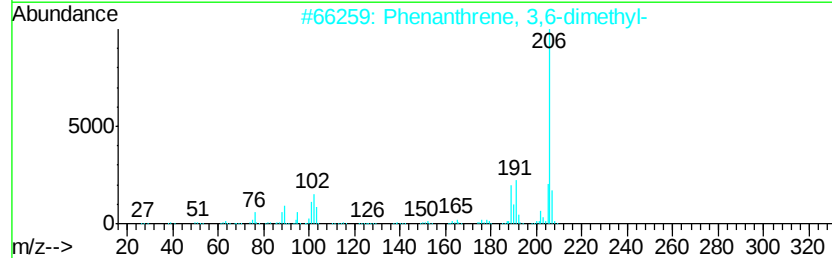
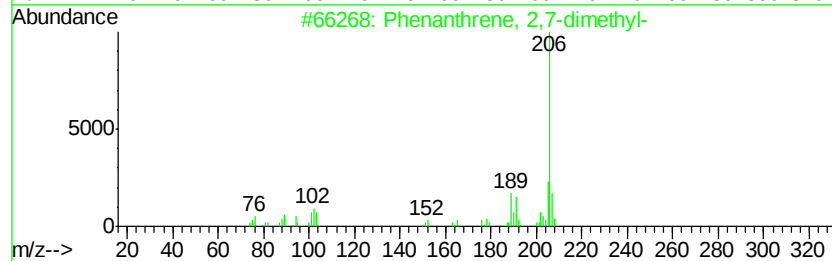
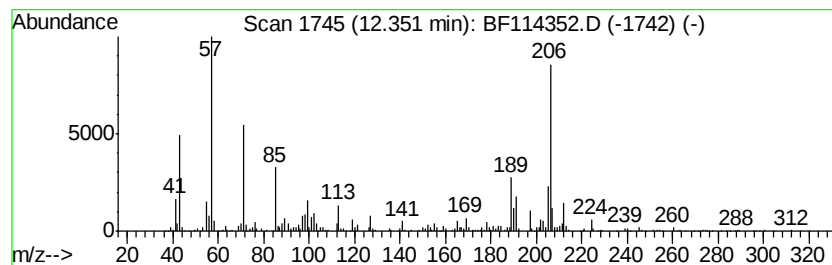
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Phenanthrene, 2,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	11.77 ng	1180000	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	64
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	64
3		1,2-Dihydro-3-phenylnaphthalene	206	C16H14	020669-52-7	55
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	50
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
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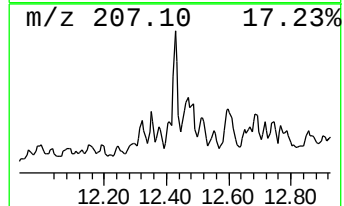
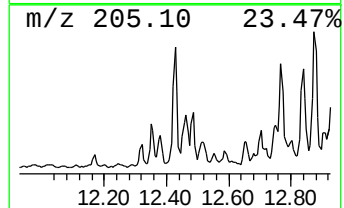
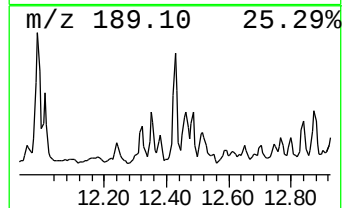
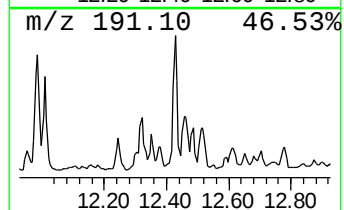
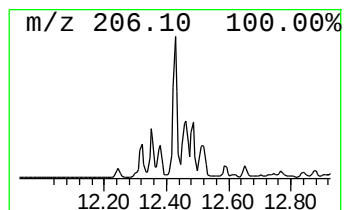
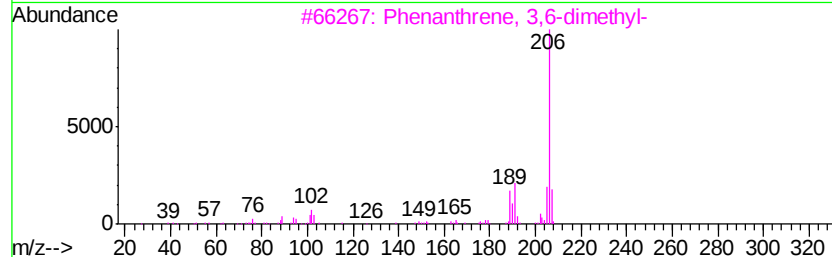
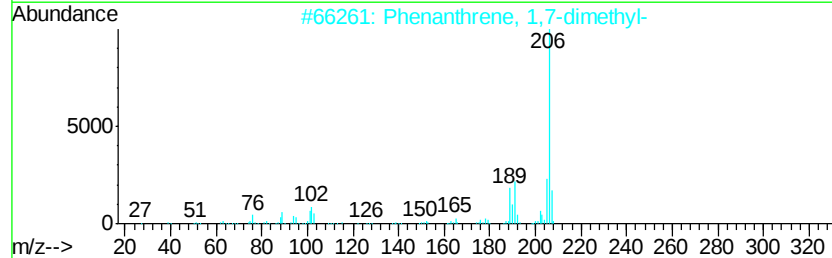
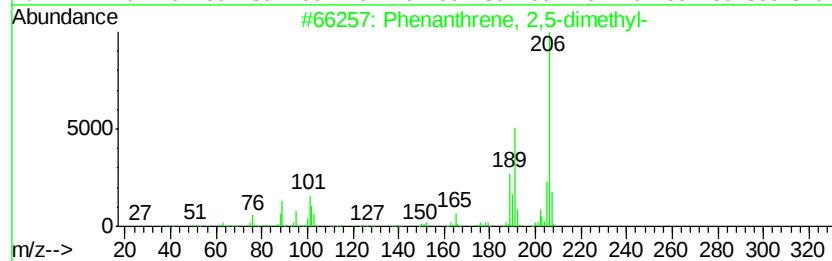
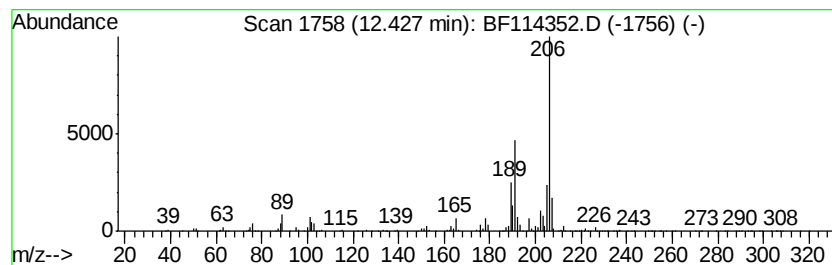
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Phenanthrene, 2,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	7.40 ng	741758	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
5		1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
Data File : BF114352.D
Acq On : 17 May 2019 16:20
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Sample : K2871-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-04(25.0-27)

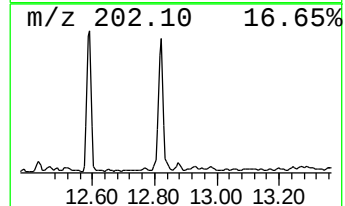
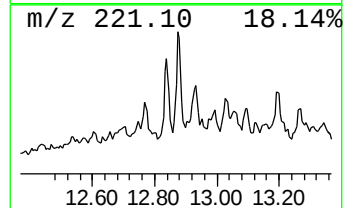
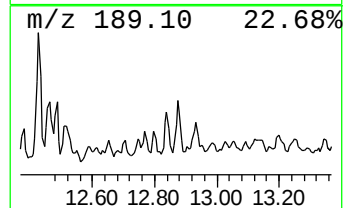
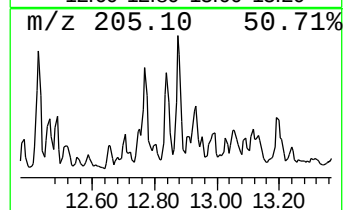
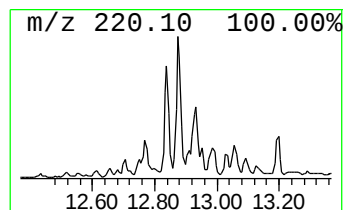
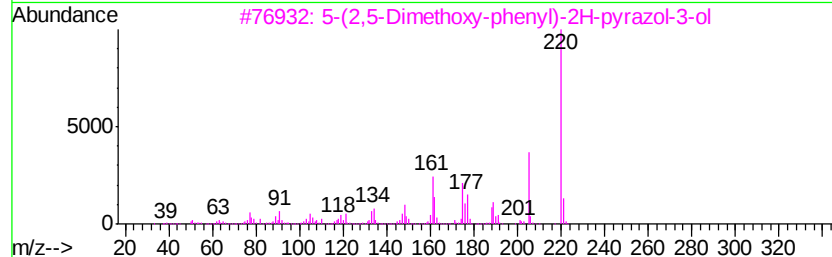
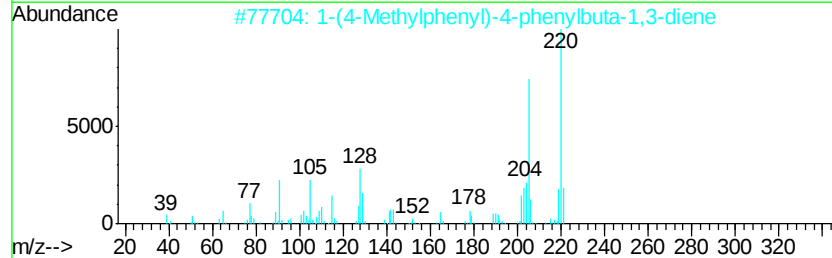
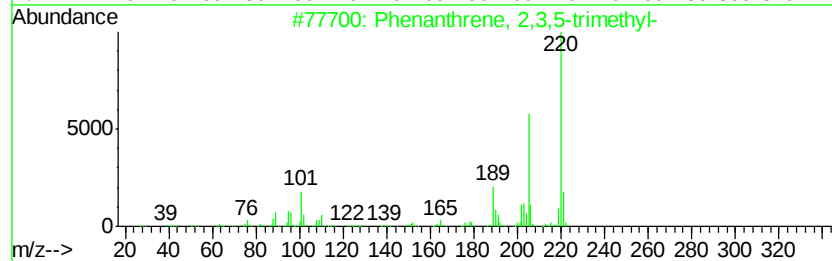
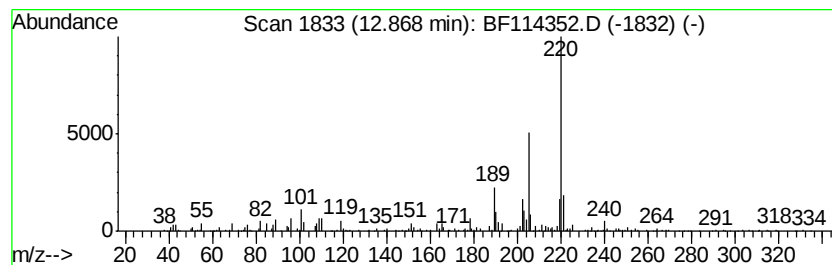
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	12.15 ng	636175	Chrysene-d12	14.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	93
2		1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	72
3		5-(2,5-Dimethoxy-phenyl)-2H-pyra...	220	C11H12N2O3	1000278-26-9	64
4		1-(Hydroxy-phenyl-methyl)-2-meth...	408	C27H36O3	108546-86-7	59
5		Phenmethylecynide, .alpha.,.alpha...	220	C11H12N2O3	1000128-94-6	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
 Data File : BF114352.D
 Acq On : 17 May 2019 16:20
 Operator : JU/SJ
 Sample : K2871-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04(25.0-27)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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
unknown6.62	6.62	55.8	ng	4162530	1	6.85	1492930	20.0
Undecane, 2,6-dim...	8.19	9.3	ng	946503	2	8.13	2036080	20.0
unknown8.42	8.42	11.4	ng	1159320	2	8.13	2036080	20.0
Nonane, 2,6-dimet...	8.55	11.0	ng	1121340	2	8.13	2036080	20.0
Dodecane, 2,7,10-...	9.15	11.4	ng	1466370	3	9.88	2561040	20.0
Dodecane, 1-chloro-	9.29	8.7	ng	1117370	3	9.88	2561040	20.0
Naphthalene, 2-(1...	9.99	6.2	ng	793791	3	9.88	2561040	20.0
Cyclohexane, (2-m...	10.14	11.3	ng	1440210	3	9.88	2561040	20.0
Naphthalene, 2,3,...	10.20	7.3	ng	934045	3	9.88	2561040	20.0
Pentadecane, 2,6,...	10.53	19.3	ng	2466440	3	9.88	2561040	20.0
Tetradecane, 4-me...	10.80	45.3	ng	4540860	4	11.37	2004510	20.0
Tridecane, 5-propyl-	11.26	35.1	ng	3518870	4	11.37	2004510	20.0
Heptadecane	11.62	11.4	ng	1139040	4	11.37	2004510	20.0
Phenanthrene, 2-m...	11.99	8.5	ng	852823	4	11.37	2004510	20.0
Phenanthrene, 2,7...	12.35	11.8	ng	1180000	4	11.37	2004510	20.0
Phenanthrene, 2,5...	12.43	7.4	ng	741758	4	11.37	2004510	20.0
Phenanthrene, 2,3...	12.87	12.2	ng	636175	5	14.03	1047570	20.0