

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111126.D
 Acq On : 6 Dec 2018 00:00
 Operator : JU/SJ
 Sample : J6201-01 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11932-72-11967

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-BF120518.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	4.369	384	388	392	rBV	96239	111117	1.68%	0.132%
2	4.999	483	495	504	rBV	712600	836583	12.68%	0.990%
3	5.410	560	565	568	rBV	6634548	5828969	88.36%	6.899%
4	6.069	674	677	683	rBV	82225	106034	1.61%	0.126%
5	6.351	722	725	727	rBV	114087	115031	1.74%	0.136%
6	6.428	732	738	743	rBV	7187074	6597167	100.00%	7.808%
7	6.569	758	762	766	rBV	6754083	6075292	92.09%	7.191%
8	6.657	772	777	782	rVB2	215149	231020	3.50%	0.273%
9	6.804	799	802	805	rBV	4111226	3239723	49.11%	3.835%
10	6.834	805	807	810	rVB	257234	188706	2.86%	0.223%
11	6.869	810	813	817	rBV3	104286	129173	1.96%	0.153%
12	6.963	825	829	832	rVB	5692553	4471023	67.77%	5.292%
13	7.093	847	851	854	rVB2	166607	211468	3.21%	0.250%
14	7.128	854	857	859	rBV2	156273	142719	2.16%	0.169%
15	7.157	859	862	866	rBV	156333	137104	2.08%	0.162%
16	7.204	867	870	873	rVV2	269314	215690	3.27%	0.255%
17	7.363	888	897	900	rBV	4271485	3980830	60.34%	4.712%
18	7.416	903	906	909	rVV	945349	724811	10.99%	0.858%
19	7.463	909	914	917	rVB4	195906	293280	4.45%	0.347%
20	7.528	921	925	928	rVV3	189025	221398	3.36%	0.262%
21	7.575	929	933	935	rVV4	85747	145100	2.20%	0.172%
22	7.598	935	937	939	rVV3	218444	223519	3.39%	0.265%
23	7.622	939	941	944	rVB2	143249	118993	1.80%	0.141%
24	7.698	949	954	955	rBV3	76221	110132	1.67%	0.130%
25	7.716	955	957	959	rVV2	148091	128393	1.95%	0.152%
26	7.734	959	960	965	rVB3	167111	201845	3.06%	0.239%
27	7.781	965	968	971	rBV2	199680	225746	3.42%	0.267%
28	7.816	971	974	978	rBV3	231321	252077	3.82%	0.298%
29	7.851	978	980	984	rVB2	333568	325595	4.94%	0.385%
30	7.893	984	987	989	rBV2	232904	217302	3.29%	0.257%
31	7.922	989	992	996	rVB3	121760	179423	2.72%	0.212%
32	7.963	996	999	1001	rBV2	87779	90363	1.37%	0.107%
33	8.034	1009	1011	1012	rBV	106175	80455	1.22%	0.095%
34	8.087	1016	1020	1023	rBV2	6401986	5480496	83.07%	6.487%

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35	8.163	1031	1033	1035	rVB	754185	502510	7.62%	0.595%
36	8.187	1035	1037	1040	rVV3	118565	101927	1.55%	0.121%
37	8.216	1040	1042	1045	rVB2	136789	103828	1.57%	0.123%
38	8.263	1045	1050	1053	rBV5	114904	233590	3.54%	0.276%
39	8.304	1055	1057	1059	rVB	190755	138413	2.10%	0.164%
40	8.392	1067	1072	1074	rBV3	297018	416394	6.31%	0.493%
41	8.410	1074	1075	1078	rVB2	168210	138627	2.10%	0.164%
42	8.445	1078	1081	1083	rBV2	279970	216540	3.28%	0.256%
43	8.475	1083	1086	1090	rVB2	392298	377361	5.72%	0.447%
44	8.522	1090	1094	1099	rBV	1066749	1091802	16.55%	1.292%
45	8.557	1099	1100	1103	rVB2	146016	113190	1.72%	0.134%
46	8.598	1104	1107	1110	rBV4	122849	116223	1.76%	0.138%
47	8.640	1111	1114	1119	rBV6	164183	188111	2.85%	0.223%
48	8.692	1119	1123	1127	rVV	1955756	1605451	24.34%	1.900%
49	8.751	1131	1133	1136	rVV3	194993	213148	3.23%	0.252%
50	8.787	1136	1139	1143	rVB3	327533	374610	5.68%	0.443%
51	8.845	1146	1149	1151	rBV3	68696	76715	1.16%	0.091%
52	8.975	1168	1171	1173	rBV3	258805	262555	3.98%	0.311%
53	9.057	1183	1185	1188	rVB	315440	236821	3.59%	0.280%
54	9.092	1188	1191	1193	rBV2	239790	221222	3.35%	0.262%
55	9.122	1193	1196	1199	rVV	997182	762361	11.56%	0.902%
56	9.163	1199	1203	1206	rVB	7495516	5619218	85.18%	6.651%
57	9.257	1215	1219	1223	rBV	1934673	1702037	25.80%	2.015%
58	9.304	1224	1227	1231	rVV5	150829	284399	4.31%	0.337%
59	9.369	1235	1238	1242	rVB6	126781	151760	2.30%	0.180%
60	9.422	1242	1247	1253	rBV3	147284	312113	4.73%	0.369%
61	9.504	1254	1261	1263	rBV4	223385	396727	6.01%	0.470%
62	9.528	1263	1265	1267	rVV2	236968	223026	3.38%	0.264%
63	9.557	1267	1270	1272	rVV	542161	639551	9.69%	0.757%
64	9.575	1272	1273	1275	rVV	819950	613389	9.30%	0.726%
65	9.598	1275	1277	1279	rVV	194834	175639	2.66%	0.208%
66	9.634	1281	1283	1285	rVB	191981	133786	2.03%	0.158%
67	9.763	1302	1305	1306	rBV2	117609	96573	1.46%	0.114%
68	9.787	1306	1309	1313	rVB	1203143	962643	14.59%	1.139%
69	9.839	1313	1318	1322	rBV2	4959274	4477275	67.87%	5.299%
70	10.022	1345	1349	1351	rBV3	99484	147651	2.24%	0.175%
71	10.045	1351	1353	1356	rVV	200664	162826	2.47%	0.193%

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 Sampling : 1 Min Area: 3 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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72	10.075	1356	1358	1361	rVV2	115826	132881	2.01%	0.157%
73	10.110	1361	1364	1367	rVB2	180417	172041	2.61%	0.204%
74	10.281	1391	1393	1396	rVB	591420	487782	7.39%	0.577%
75	10.339	1401	1403	1408	rVB5	76049	89890	1.36%	0.106%
76	10.504	1427	1431	1434	rBV	286181	269646	4.09%	0.319%
77	10.622	1447	1451	1454	rBV	3612774	2865398	43.43%	3.392%
78	10.757	1471	1474	1475	rBV	336397	293748	4.45%	0.348%
79	10.769	1475	1476	1481	rVB	282919	230208	3.49%	0.272%
80	11.204	1547	1550	1552	rBV	190090	150437	2.28%	0.178%
81	11.233	1552	1555	1559	rVB	130220	120361	1.82%	0.142%
82	11.322	1566	1570	1573	rBV	4342878	3498447	53.03%	4.141%
83	11.628	1619	1622	1625	rBV	125465	111676	1.69%	0.132%
84	11.857	1657	1661	1671	rBV2	564969	640827	9.71%	0.758%
85	12.528	1773	1775	1780	rBV	110606	110198	1.67%	0.130%
86	12.639	1791	1794	1798	rBV	303256	311461	4.72%	0.369%
87	12.904	1835	1839	1843	rBV	4616354	3934191	59.63%	4.657%
88	13.816	1990	1994	1998	rBV3	234495	329497	4.99%	0.390%
89	13.951	2013	2017	2021	rBV	2941497	2878347	43.63%	3.407%
90	14.445	2098	2101	2104	rBV	139966	161728	2.45%	0.191%
91	14.816	2162	2164	2168	rVB	395546	381454	5.78%	0.451%
92	15.074	2206	2208	2211	rVB	191148	178576	2.71%	0.211%
93	15.392	2257	2262	2266	rBV	1604552	1911749	28.98%	2.263%

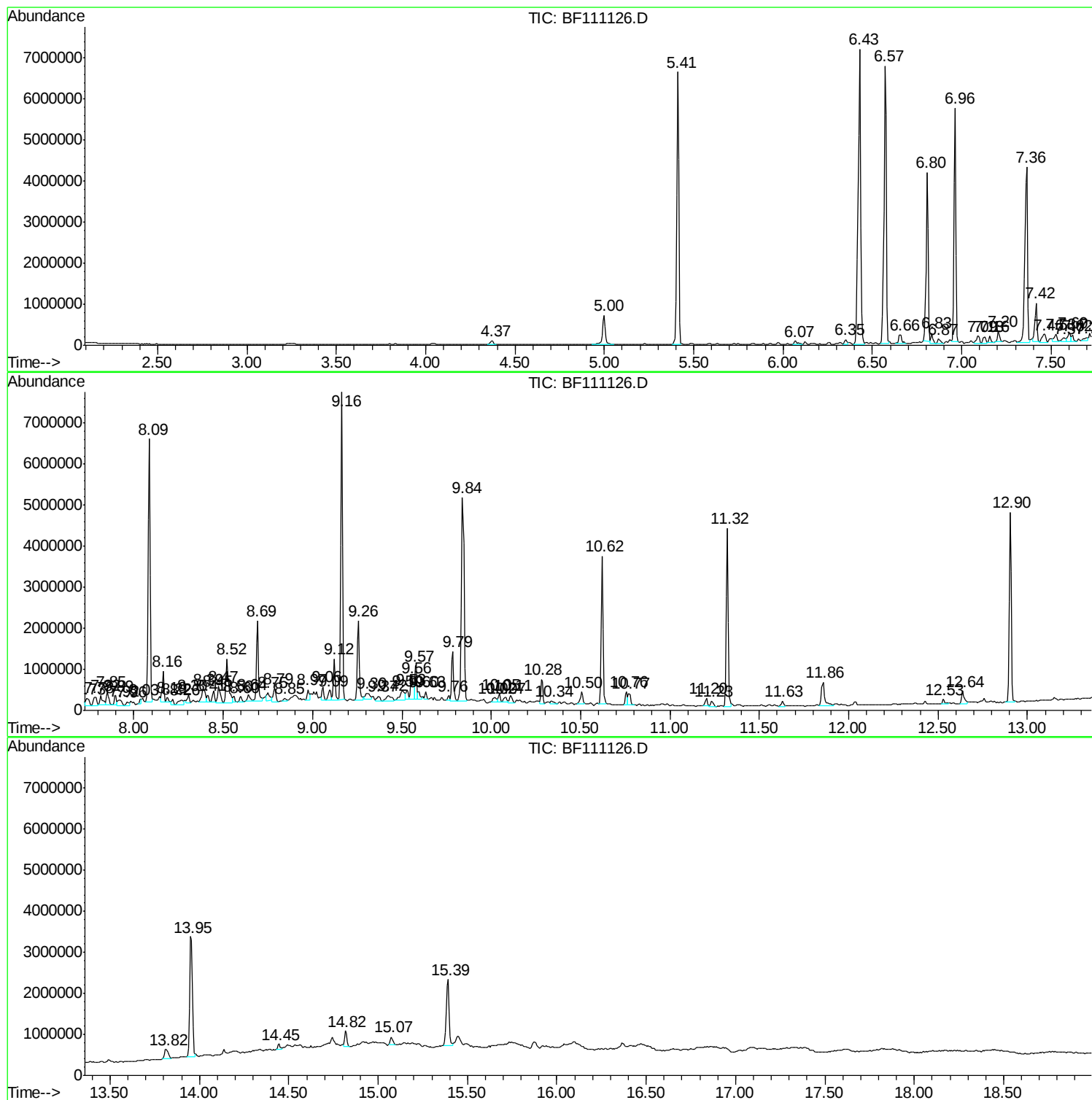
Sum of corrected areas: 84487132

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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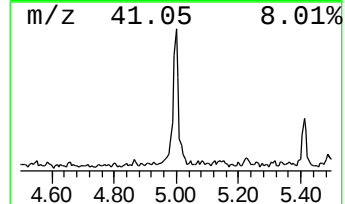
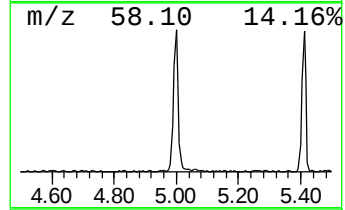
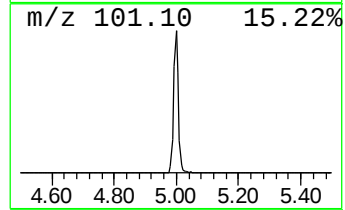
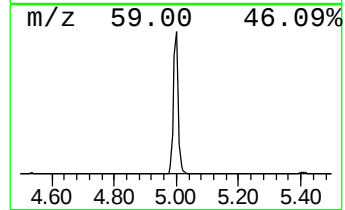
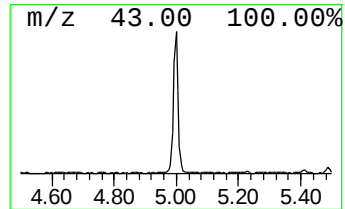
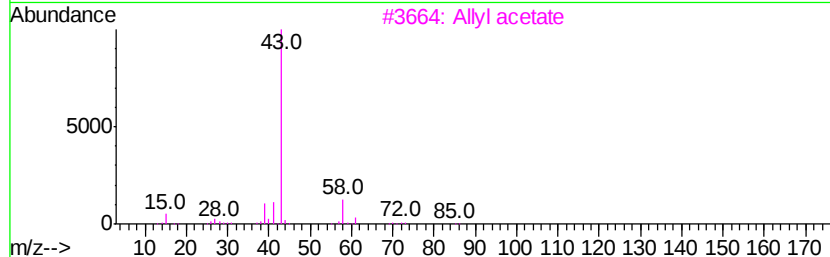
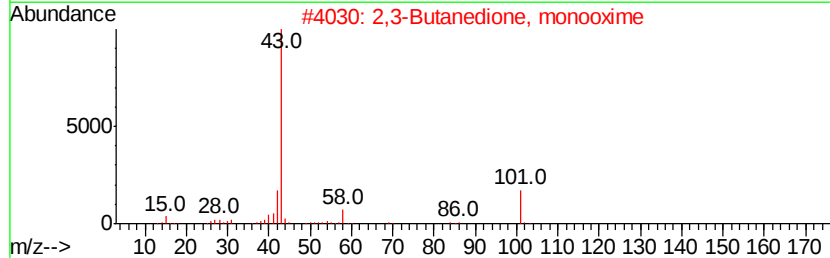
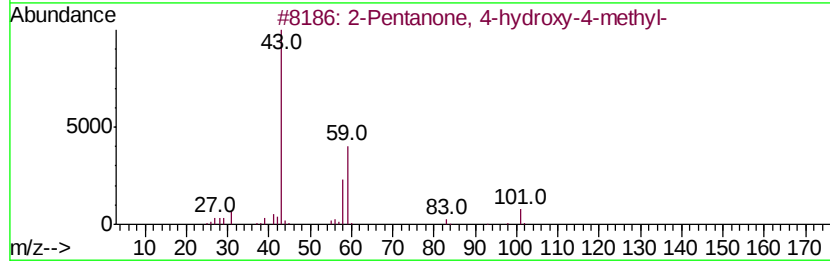
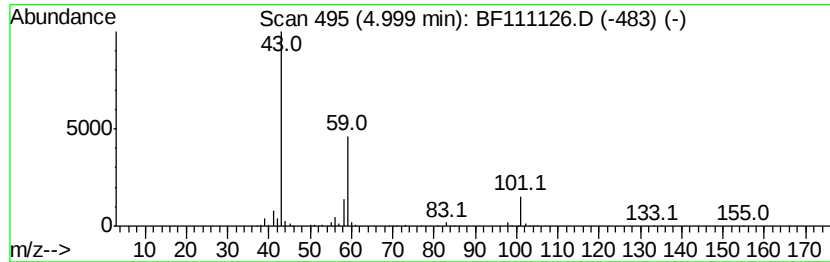
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.00	5.16 ng	836583	1,4-Dichlorobenzene-d4	6.80	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
3	Allyl acetate	100	C5H8O2	000591-87-7	9
4	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	7
5	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	7



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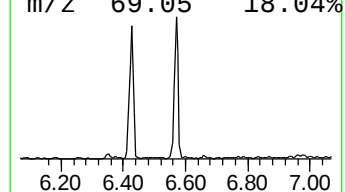
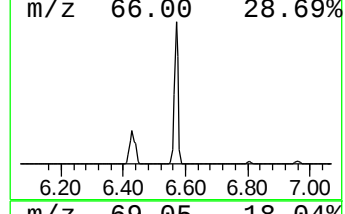
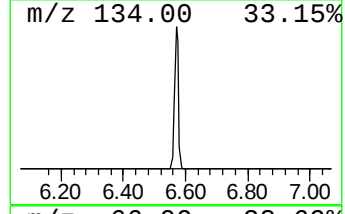
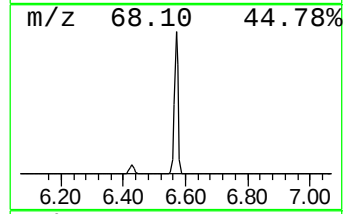
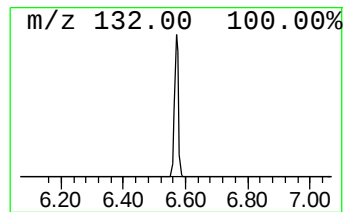
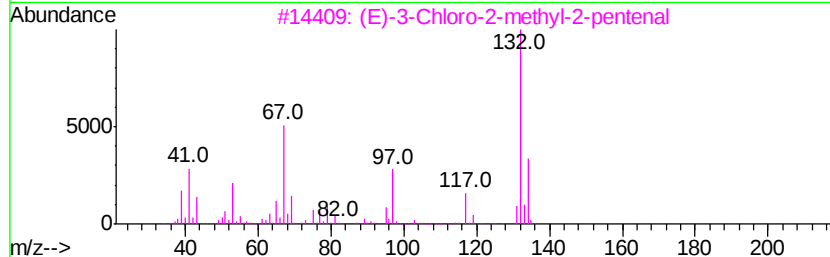
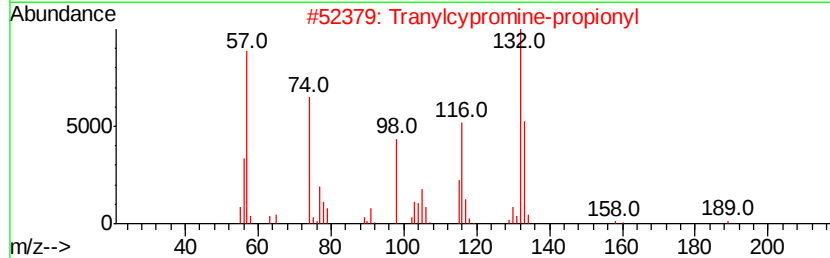
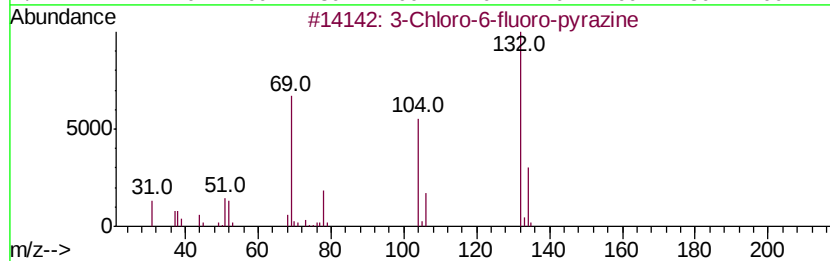
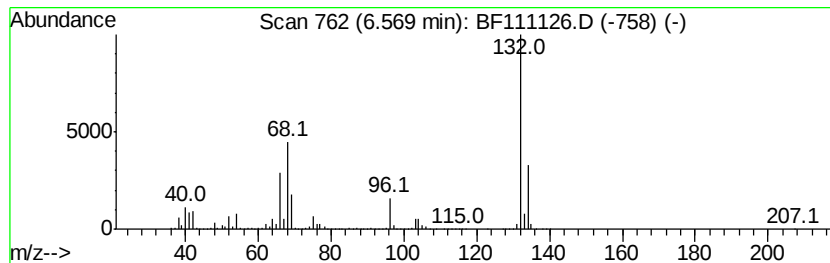
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	37.51 ng	6075290	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
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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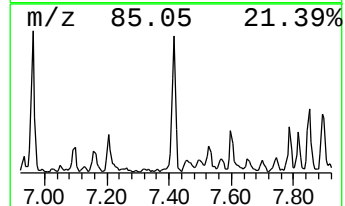
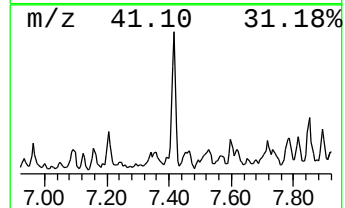
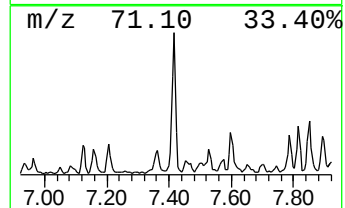
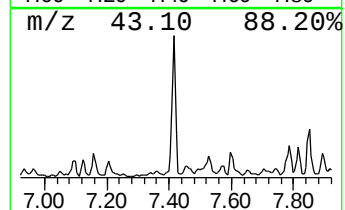
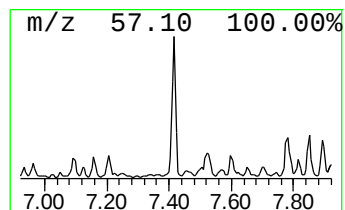
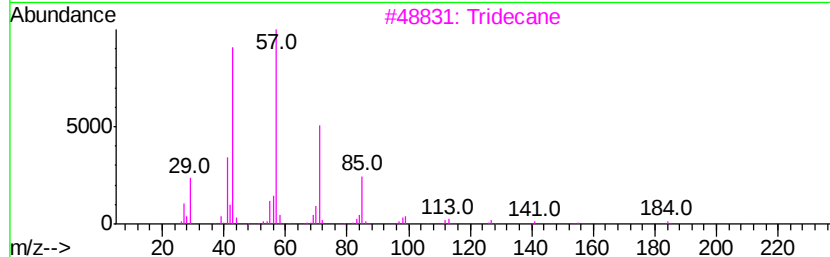
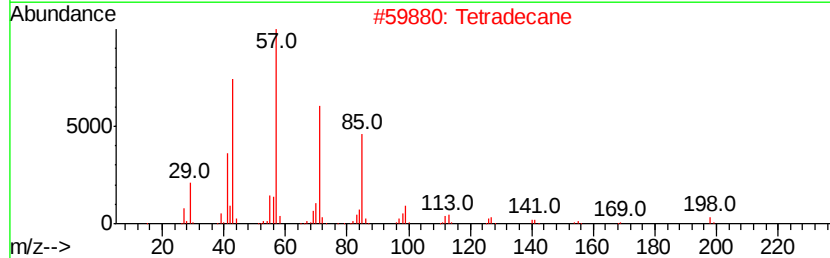
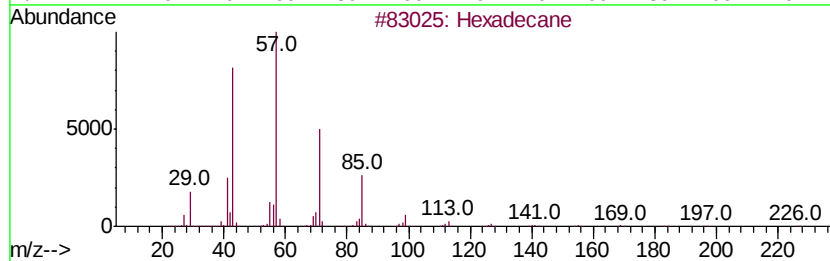
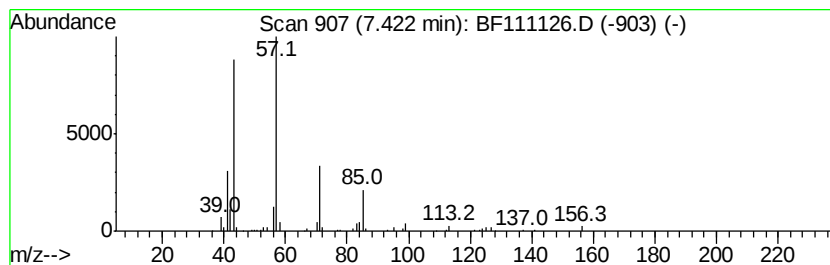
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	4.47 ng	724811	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	83
2		Tetradecane	198	C14H30	000629-59-4	83
3		Tridecane	184	C13H28	000629-50-5	83
4		Pentadecane	212	C15H32	000629-62-9	72
5		1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
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Acq On : 6 Dec 2018 00:00
Operator : JU/SJ
Sample : J6201-01 2X
Misc :
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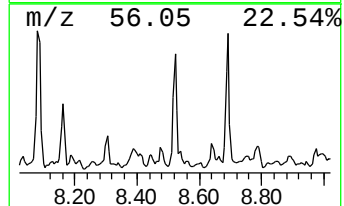
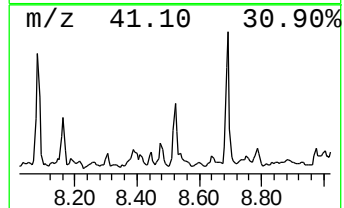
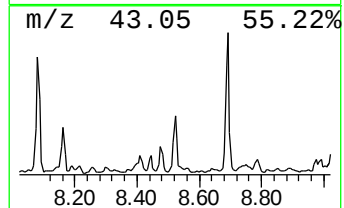
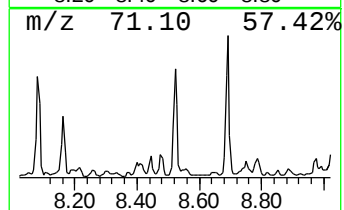
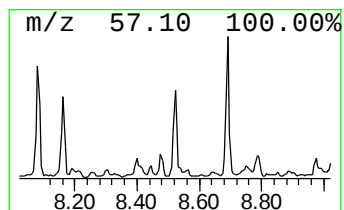
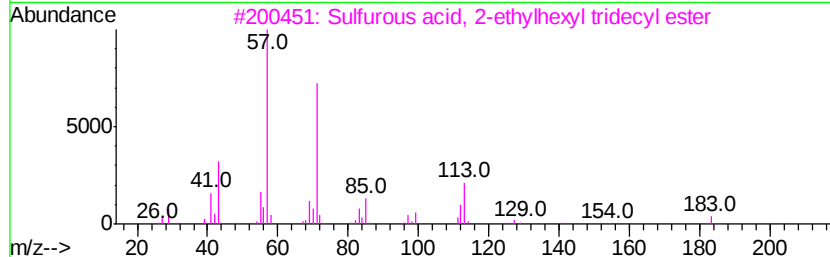
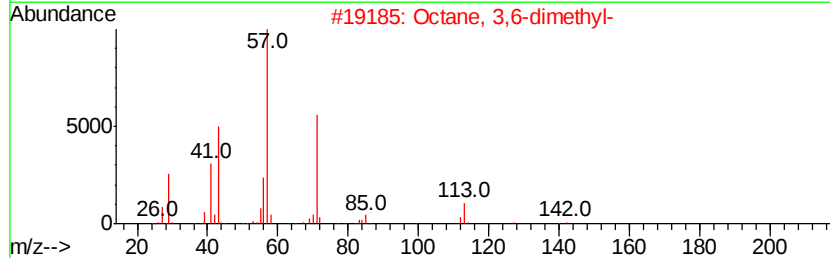
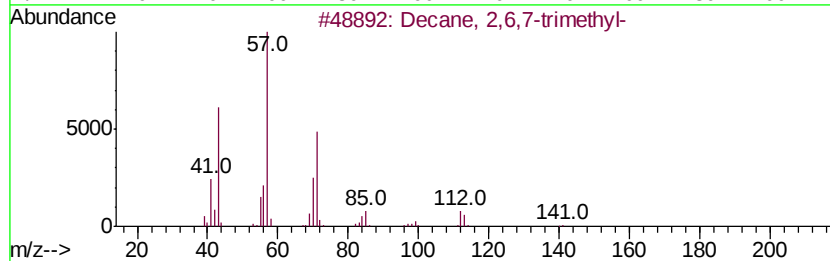
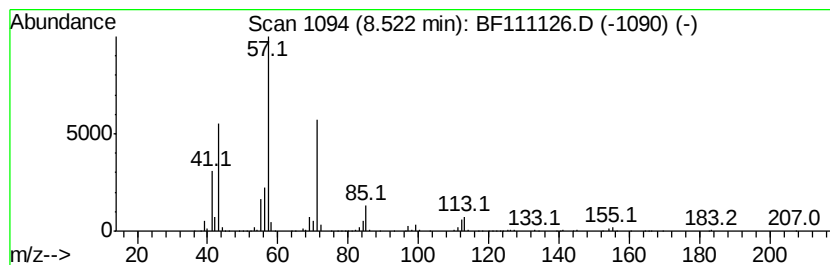
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Decane, 2,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.52	3.98 ng	1091800	Naphthalene-d8	8.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 2,6,7-trimethyl-	184 C13H28	062108-25-2	83
2	Octane, 3,6-dimethyl-	142 C10H22	015869-94-0	64
3	Sulfurous acid, 2-ethylhexyl tri...	376 C21H44O3S	1000309-19-6	64
4	Heptadecane, 2,6-dimethyl-	268 C19H40	054105-67-8	64
5	Nonane, 3-methyl-	142 C10H22	005911-04-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
Data File : BF111126.D
Acq On : 6 Dec 2018 00:00
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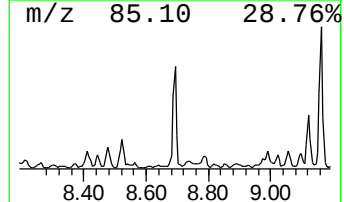
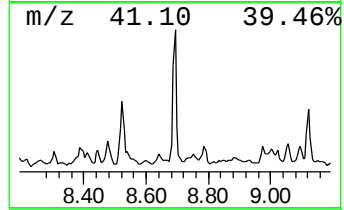
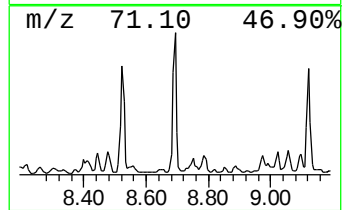
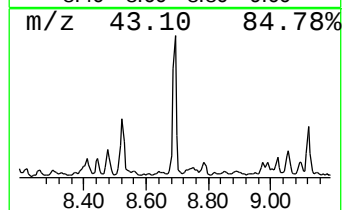
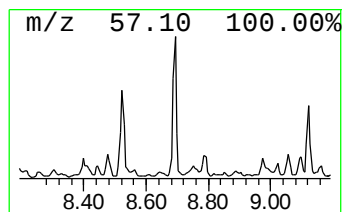
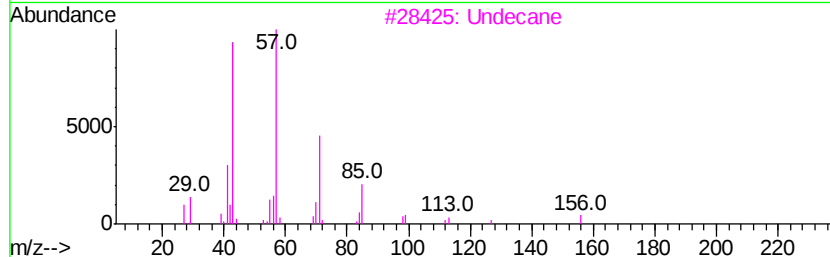
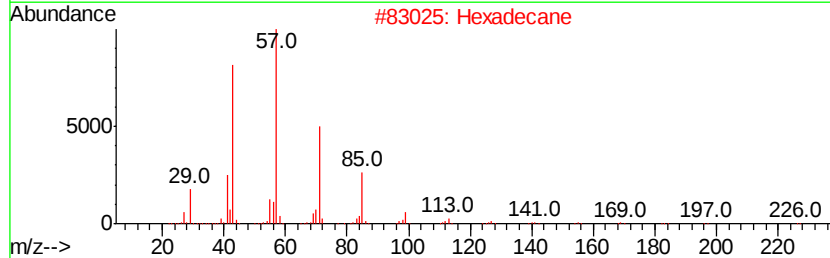
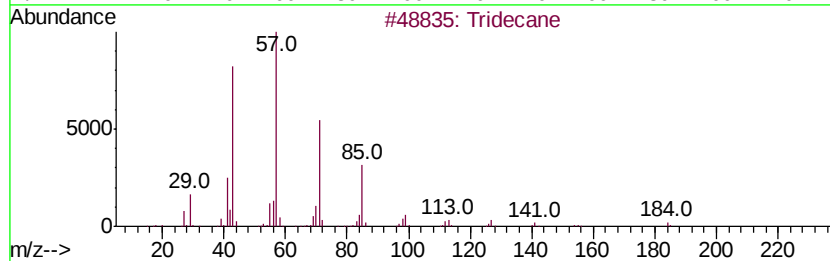
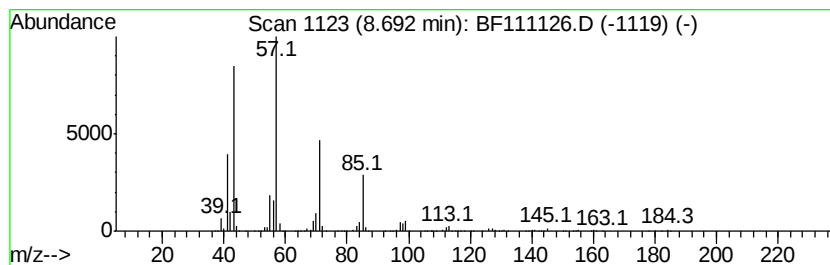
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.69	5.86 ng	1605450	Naphthalene-d8	8.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	96
2	Hexadecane	226	C16H34	000544-76-3	90
3	Undecane	156	C11H24	001120-21-4	86
4	Dodecane	170	C12H26	000112-40-3	72
5	Nonane, 5-butyl-	184	C13H28	017312-63-9	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
Data File : BF111126.D
Acq On : 6 Dec 2018 00:00
Operator : JU/SJ
Sample : J6201-01 2X
Misc :
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Instrument :
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ClientSampleId :
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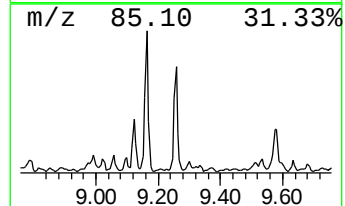
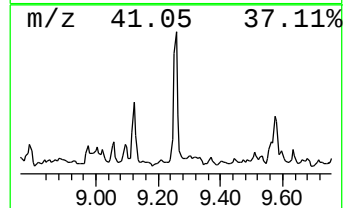
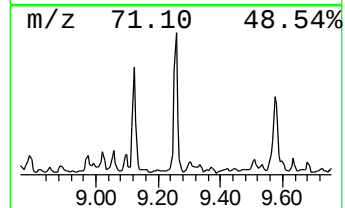
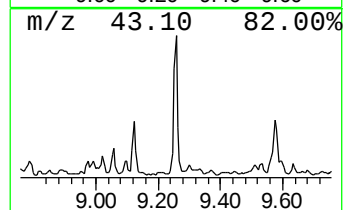
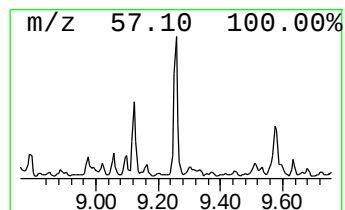
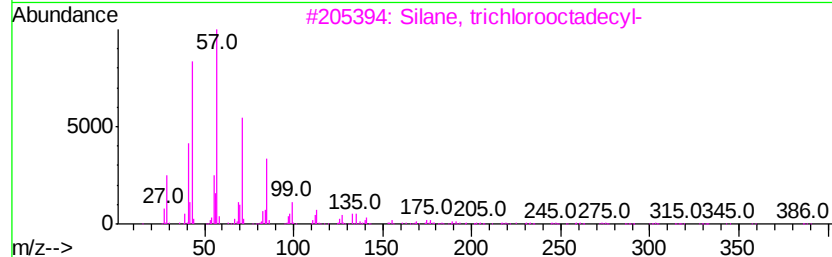
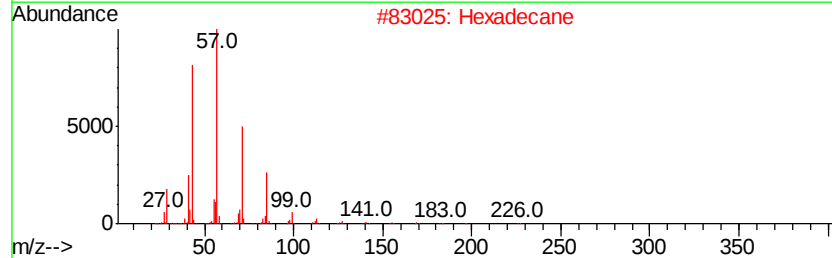
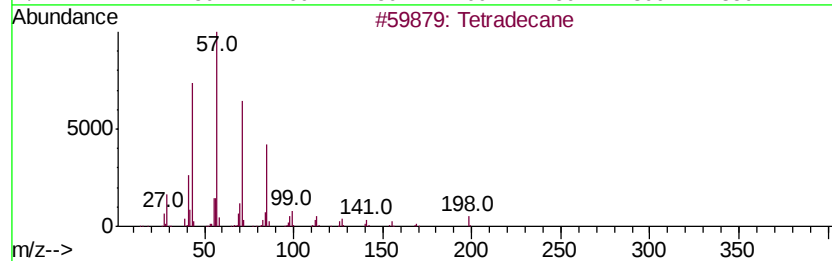
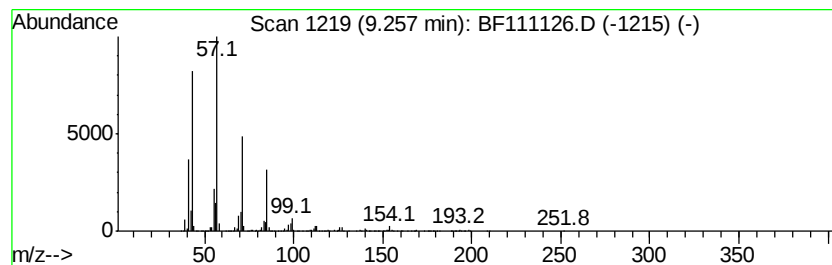
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.26	7.60 ng	1702040	Acenaphthene-d10	9.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	94
2		Hexadecane	226	C16H34	000544-76-3	91
3		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	86
4		Undecane	156	C11H24	001120-21-4	86
5		Tridecane, 7-propyl-	226	C16H34	055045-09-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
Data File : BF111126.D
Acq On : 6 Dec 2018 00:00
Operator : JU/SJ
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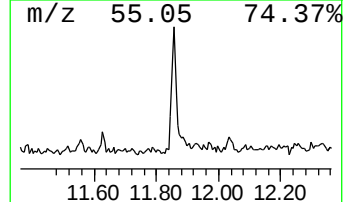
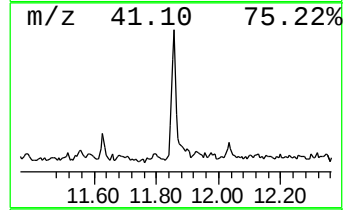
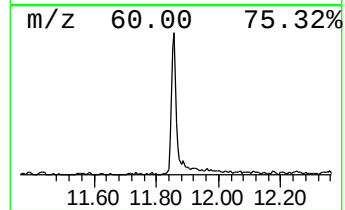
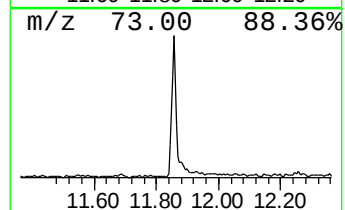
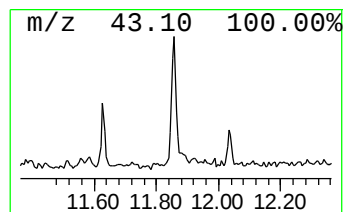
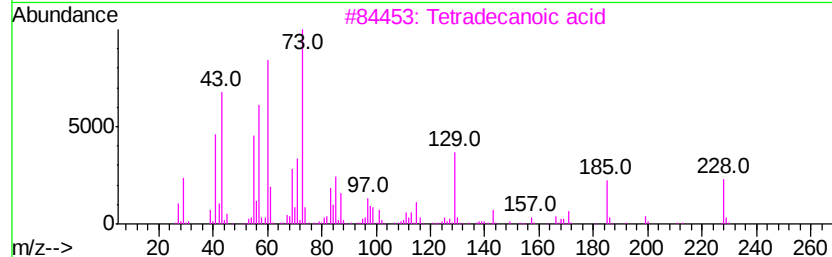
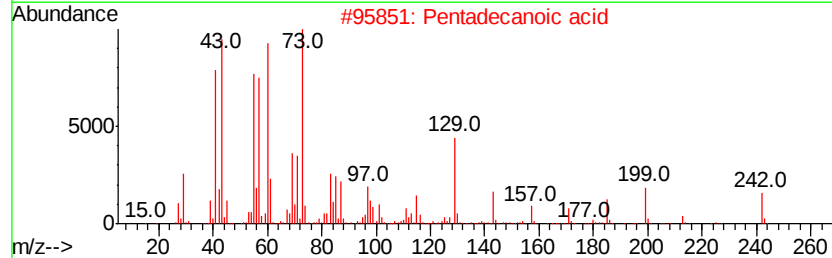
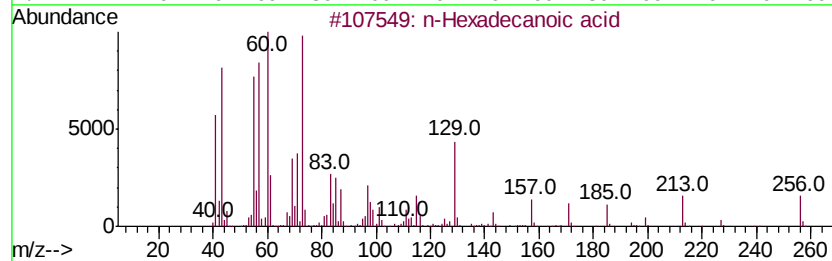
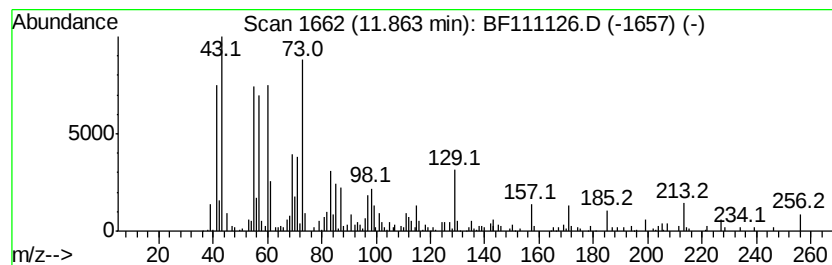
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	3.66 ng	640827	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
4			Tridecanoic acid	214	C13H26O2	000638-53-9	74
5			Dodecanoic acid	200	C12H24O2	000143-07-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
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Sample : J6201-01 2X
Misc :
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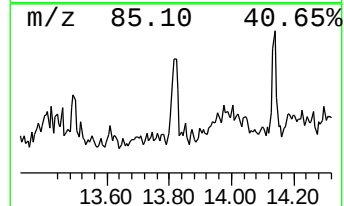
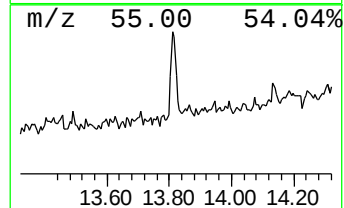
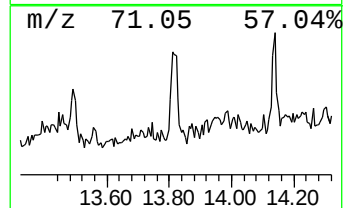
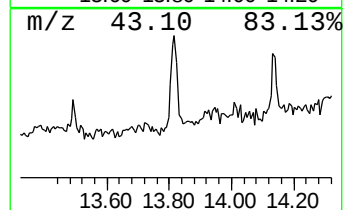
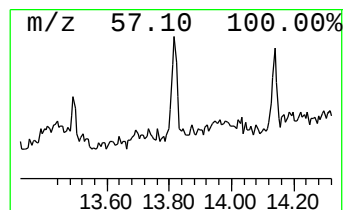
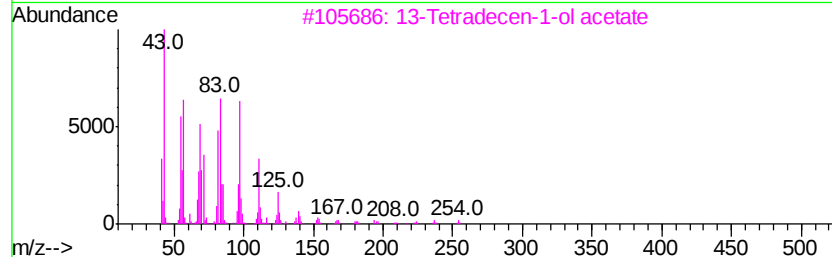
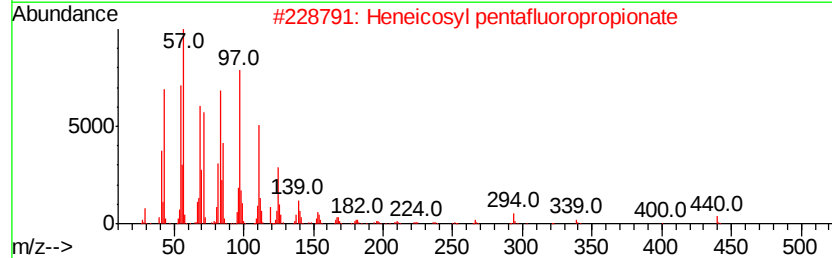
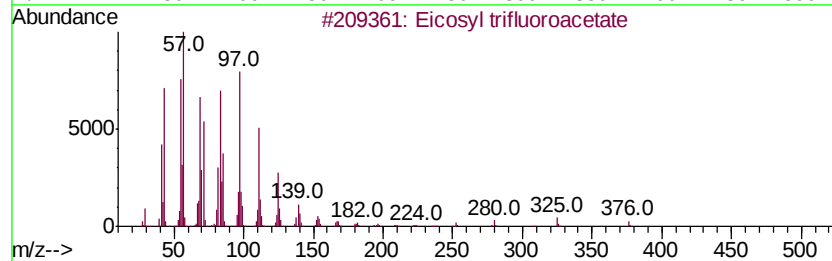
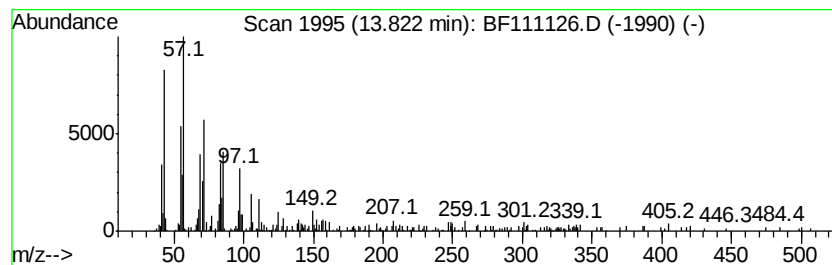
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosyl trifluoroacetate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.82	2.29 ng	329497	Chrysene-d12	13.96	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	90
2	Heneicosyl pentafluoropropionate	458	C24H43F5O2	1000351-81-1	90
3	13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	89
4	Docosyl trifluoroacetate	422	C24H45F3O2	1000351-75-5	87
5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
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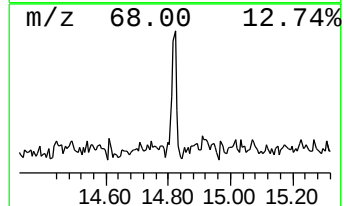
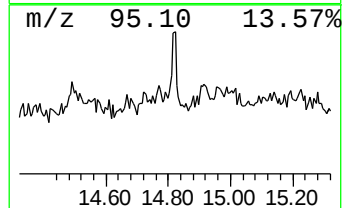
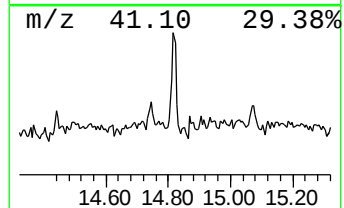
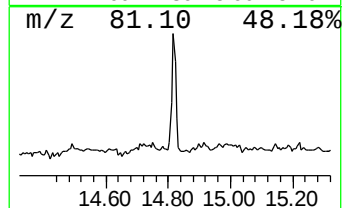
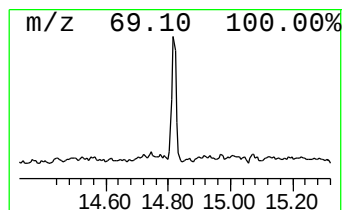
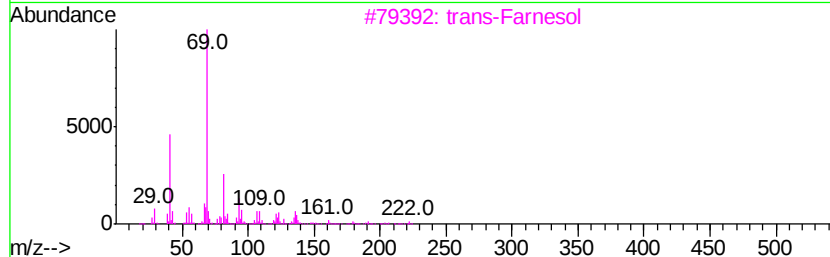
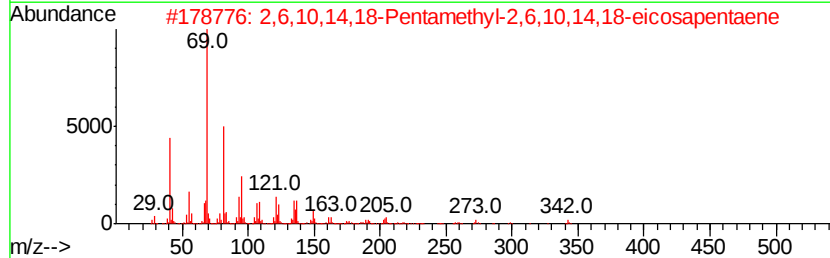
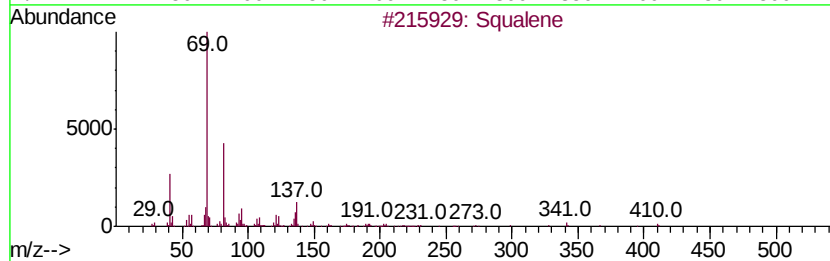
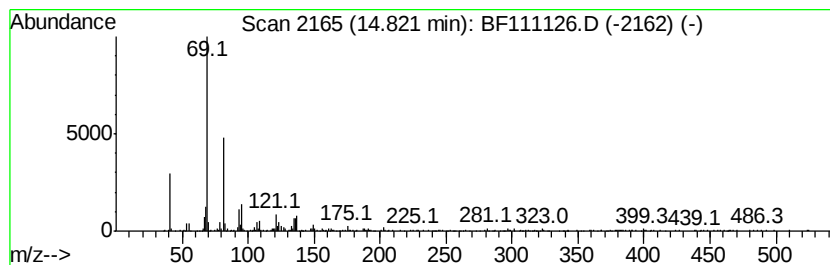
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	3.99 ng	381454	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	80
2		2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	74
3		trans-Farnesol	222	C15H26O	000106-28-5	64
4		2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	53
5		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120518\
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Sample : J6201-01 2X
Misc :
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ClientSampleId :
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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
2-Pentanone, 4-hy...	5.00	5.2 ng		836583	1	6.80	3239720	20.0
unknown6.57	6.57	37.5 ng		6075290	1	6.80	3239720	20.0
Hexadecane	7.42	4.5 ng		724811	1	6.80	3239720	20.0
Decane, 2,6,7-tri...	8.52	4.0 ng		1091800	2	8.09	5480500	20.0
Tridecane	8.69	5.9 ng		1605450	2	8.09	5480500	20.0
Tetradecane	9.26	7.6 ng		1702040	3	9.84	4477280	20.0
n-Hexadecanoic acid	11.86	3.7 ng		640827	4	11.32	3498450	20.0
Eicosyl trifluoro...	13.82	2.3 ng		329497	5	13.96	2878350	20.0
Squalene	14.82	4.0 ng		381454	6	15.39	1911750	20.0