

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106700.D
 Acq On : 22 Jun 2018 11:18
 Operator : JU/SJ
 Sample : J3603-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26KV-TP-10

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-BF061918.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	5.201	483	487	501	rBB	122237	161151	9.46%	0.990%
2	5.607	552	556	559	rBV	1418368	1391295	81.66%	8.543%
3	6.607	721	726	729	rBV	1377249	1500890	88.10%	9.216%
4	6.736	744	748	752	rBV	1453627	1510226	88.64%	9.273%
5	6.778	752	755	759	rVB	27725	23537	1.38%	0.145%
6	6.960	782	786	789	rBV	510803	407605	23.92%	2.503%
7	7.119	808	813	816	rVB	1364186	1236410	72.57%	7.592%
8	7.525	877	882	885	rBV	1032117	1104200	64.81%	6.780%
9	7.601	889	895	898	rBV2	14353	17696	1.04%	0.109%
10	8.242	1000	1004	1007	rBV	623965	524929	30.81%	3.223%
11	8.642	1069	1072	1075	rBV	22091	20514	1.20%	0.126%
12	9.319	1180	1187	1190	rBV	1761881	1703679	100.00%	10.461%
13	9.383	1195	1198	1201	rBV2	21115	19566	1.15%	0.120%
14	9.707	1249	1253	1260	rVB	271791	229785	13.49%	1.411%
15	9.913	1284	1288	1292	rBV	36854	32649	1.92%	0.200%
16	10.001	1299	1303	1307	rVB2	694013	604275	35.47%	3.710%
17	10.413	1370	1373	1376	rVB	57825	49680	2.92%	0.305%
18	10.630	1405	1410	1412	rBV	20615	21175	1.24%	0.130%
19	10.795	1434	1438	1442	rBV	1197870	1179737	69.25%	7.244%
20	10.883	1450	1453	1459	rBV2	60177	76743	4.50%	0.471%
21	11.336	1526	1530	1532	rBV2	45247	40925	2.40%	0.251%
22	11.495	1553	1557	1559	rBV	789520	665694	39.07%	4.088%
23	11.519	1559	1561	1564	rVB	53613	43294	2.54%	0.266%
24	11.760	1600	1602	1605	rVB2	41292	33381	1.96%	0.205%
25	11.824	1610	1613	1618	rVB5	12360	18832	1.11%	0.116%
26	11.995	1639	1642	1644	rBV	30126	26064	1.53%	0.160%
27	12.024	1644	1647	1651	rVB3	44122	45838	2.69%	0.281%
28	12.101	1658	1660	1663	rBV2	32812	27211	1.60%	0.167%
29	12.171	1668	1672	1675	rBV2	31602	30425	1.79%	0.187%
30	12.289	1688	1692	1695	rBV2	26757	24862	1.46%	0.153%
31	12.542	1732	1735	1740	rBV2	50736	68393	4.01%	0.420%
32	12.630	1747	1750	1754	rVB2	16849	20377	1.20%	0.125%
33	12.707	1760	1763	1767	rVB	86307	76427	4.49%	0.469%
34	12.942	1797	1803	1808	rBV3	66713	99368	5.83%	0.610%

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.989	1808	1811	1814	rVB2	24735	23904	1.40%	0.147%
36	13.077	1822	1826	1830	rBV	1618352	1659536	97.41%	10.190%
37	13.154	1835	1839	1846	rVB3	19243	34129	2.00%	0.210%
38	13.777	1943	1945	1950	rVB2	19715	23271	1.37%	0.143%
39	13.960	1972	1976	1979	rBV2	69099	68821	4.04%	0.423%
40	14.077	1989	1996	1998	rBV4	15746	25483	1.50%	0.156%
41	14.148	2003	2008	2011	rVV	629032	594224	34.88%	3.649%
42	14.171	2011	2012	2015	rVB	40925	24362	1.43%	0.150%
43	14.289	2029	2032	2035	rBV3	13868	18757	1.10%	0.115%
44	14.548	2070	2076	2079	rBV	20348	34948	2.05%	0.215%
45	14.607	2083	2086	2092	rVB6	16106	24522	1.44%	0.151%
46	15.236	2189	2193	2196	rBV	30092	44904	2.64%	0.276%
47	15.554	2244	2247	2252	rVB	25250	32701	1.92%	0.201%
48	15.624	2255	2259	2265	rVB	28129	40074	2.35%	0.246%
49	15.695	2265	2271	2280	rBV	420459	581637	34.14%	3.571%
50	16.136	2344	2346	2352	rVB6	10841	17911	1.05%	0.110%

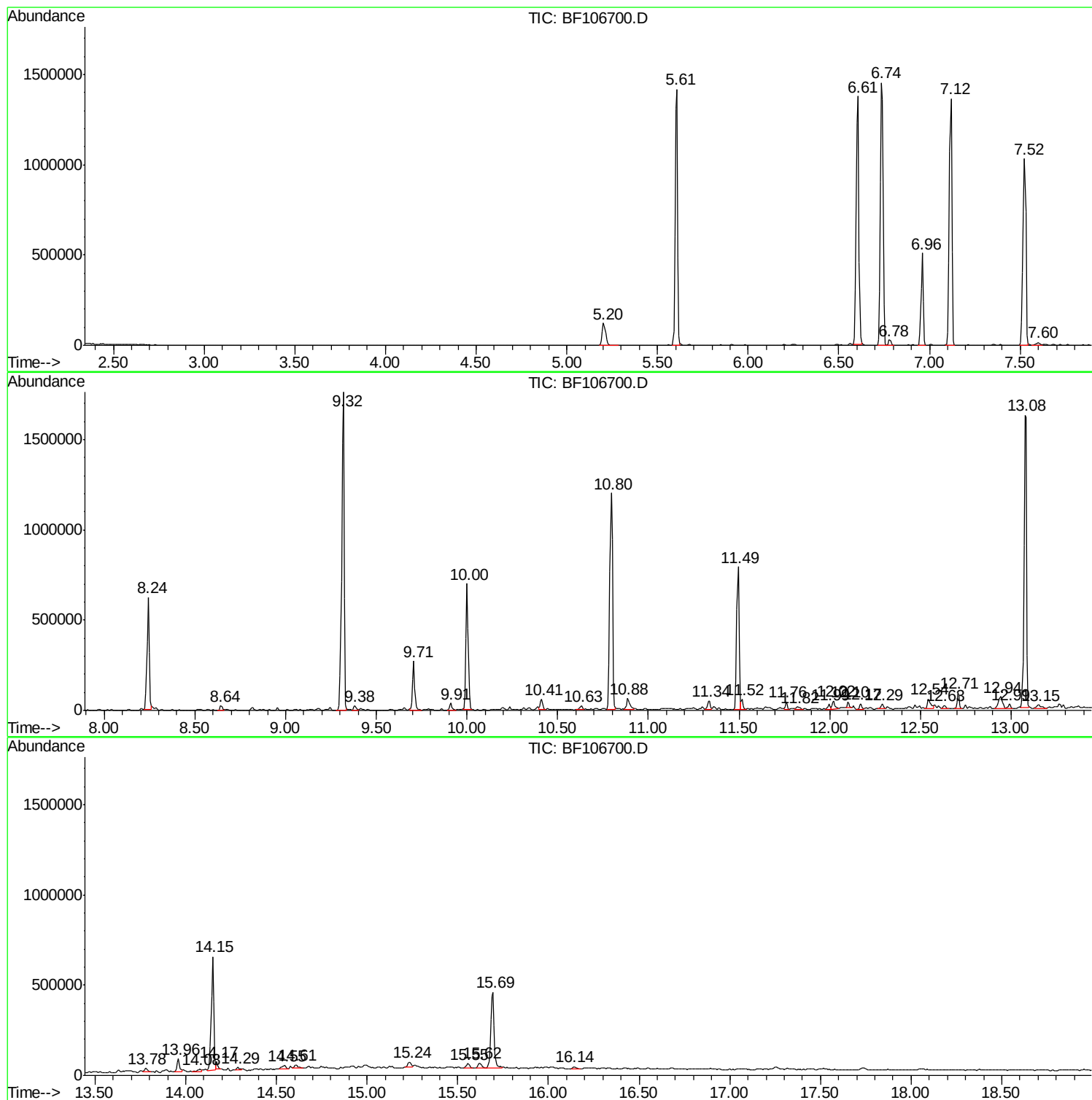
Sum of corrected areas: 16286017

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
26KV-TP-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

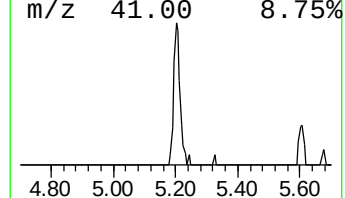
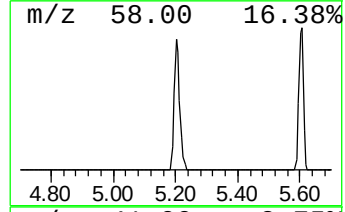
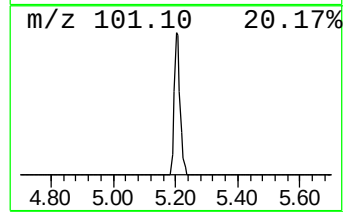
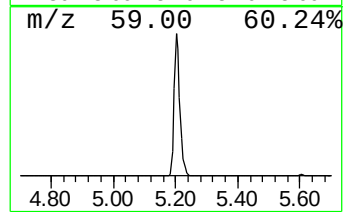
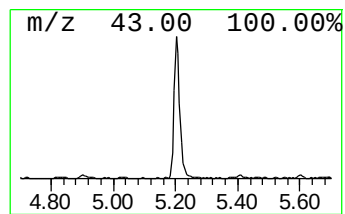
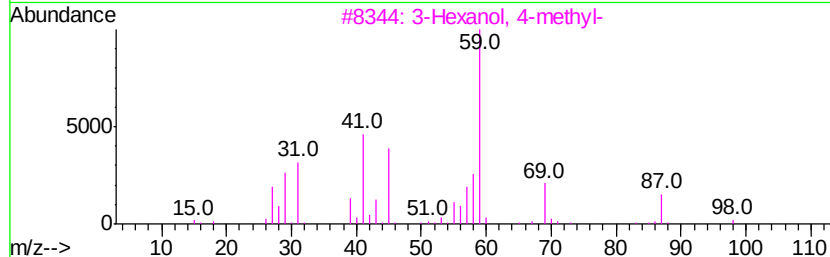
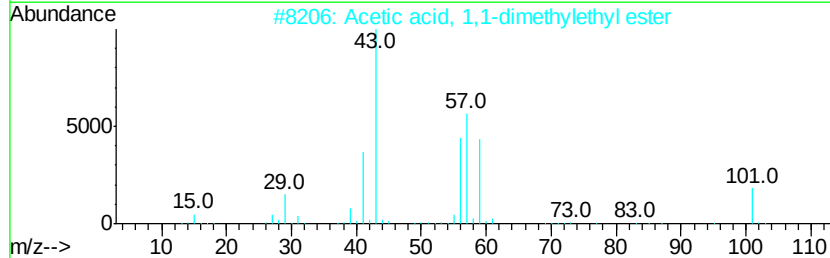
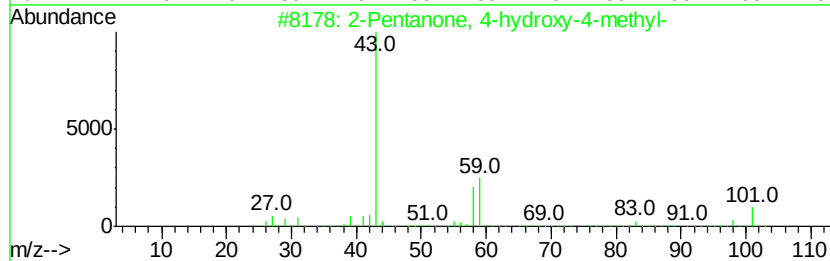
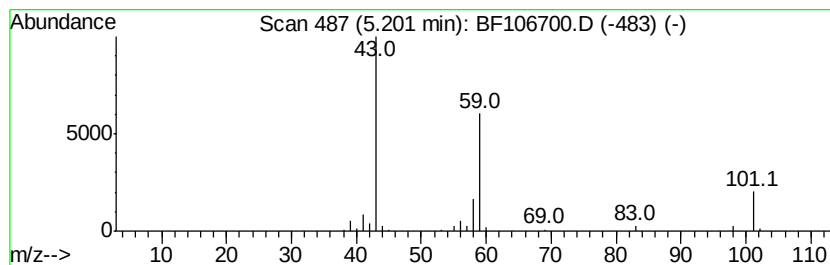
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	7.91 ng	161151	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

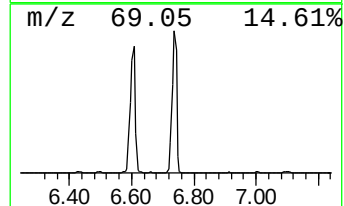
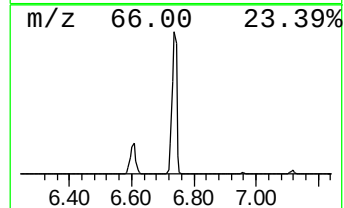
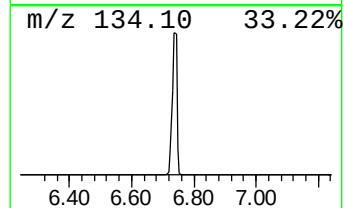
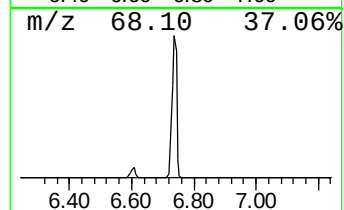
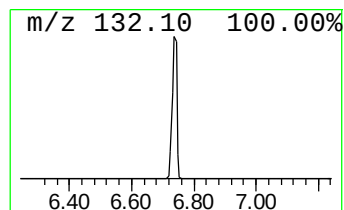
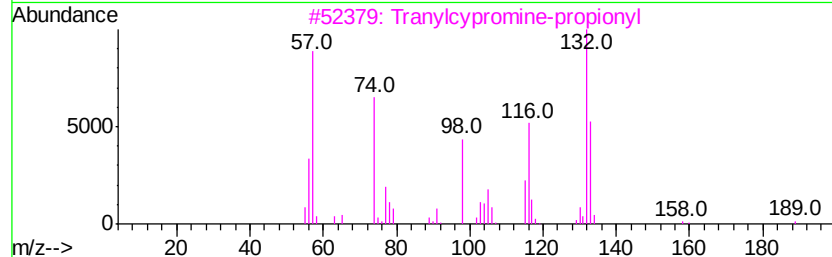
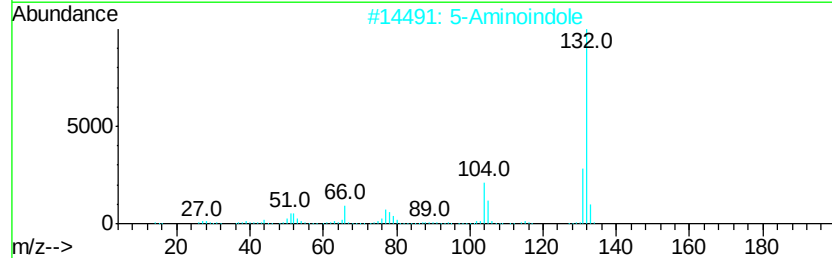
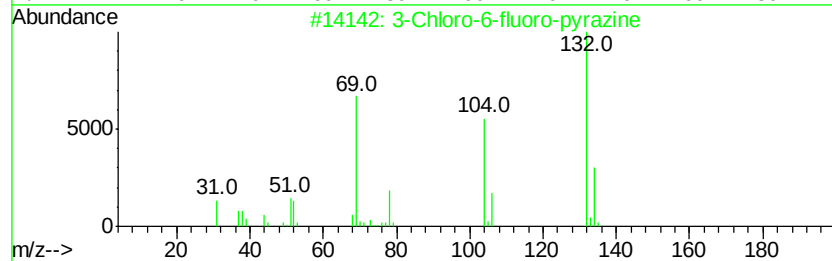
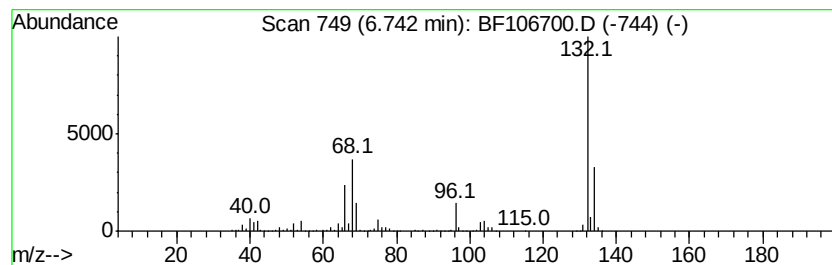
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	74.10 ng	1510230	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

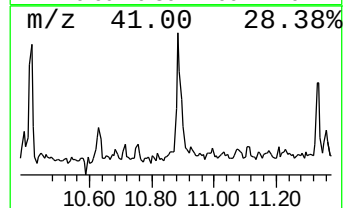
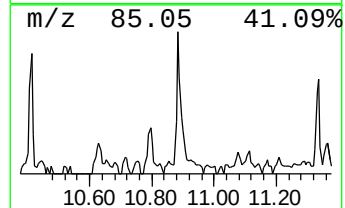
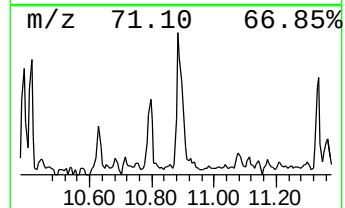
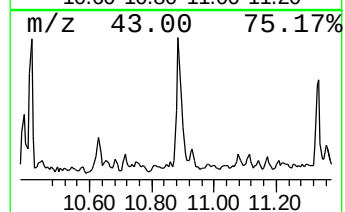
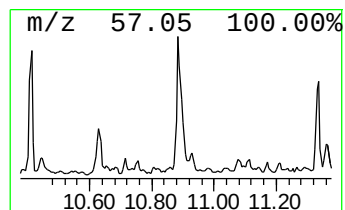
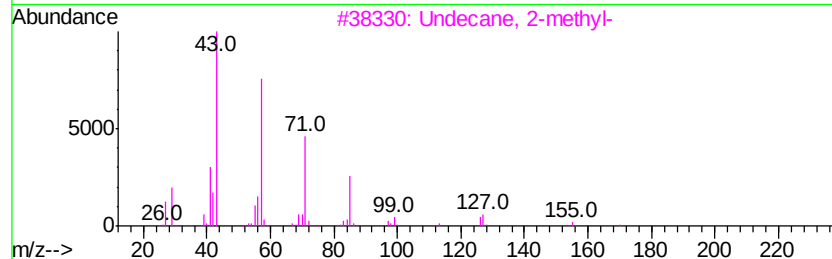
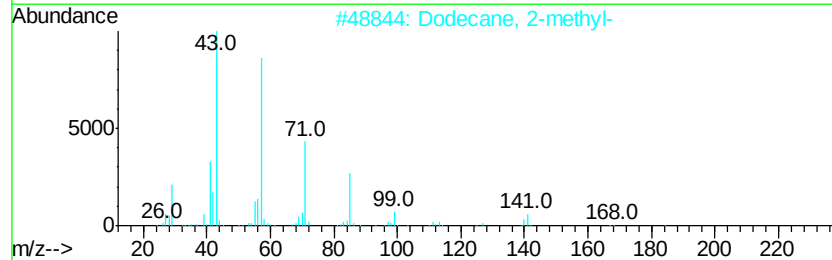
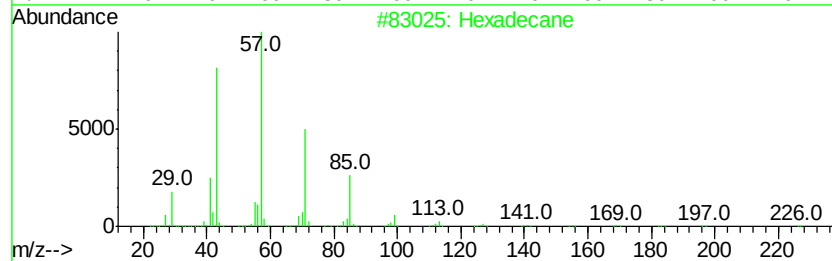
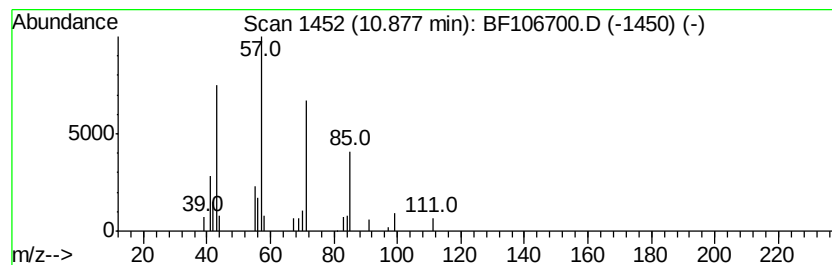
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
10.88	2.31 ng	76743	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	80
2	Dodecane, 2-methyl-		184	C13H28	001560-97-0	80
3	Undecane, 2-methyl-		170	C12H26	007045-71-8	72
4	Heneicosane		296	C21H44	000629-94-7	72
5	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

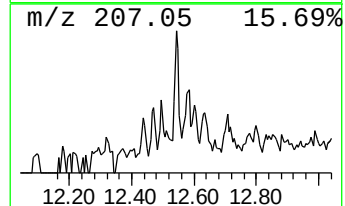
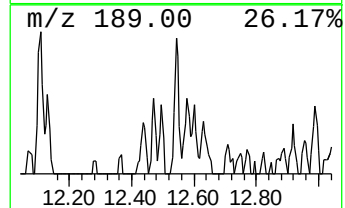
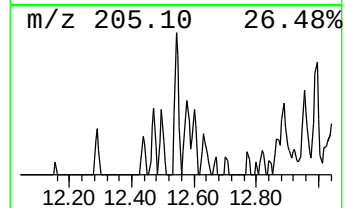
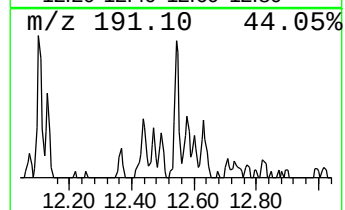
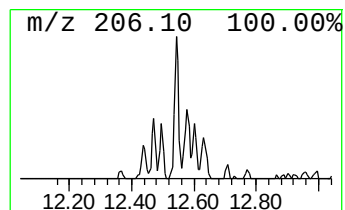
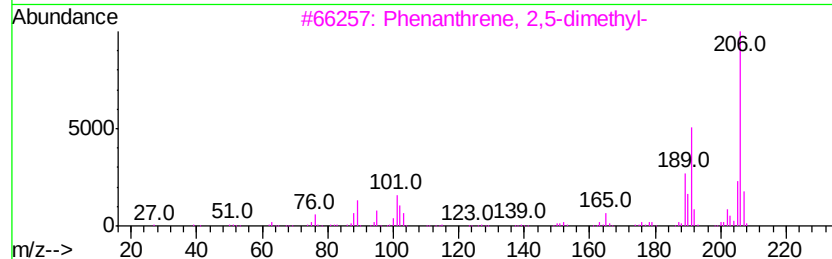
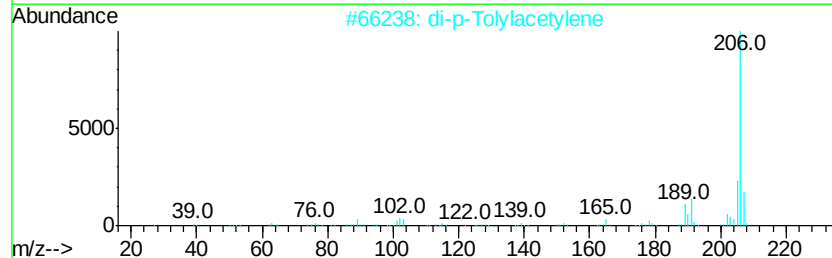
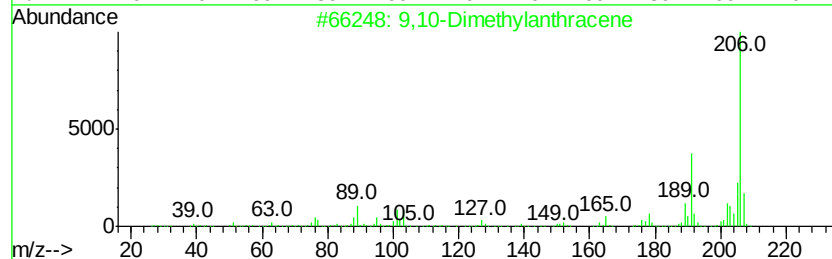
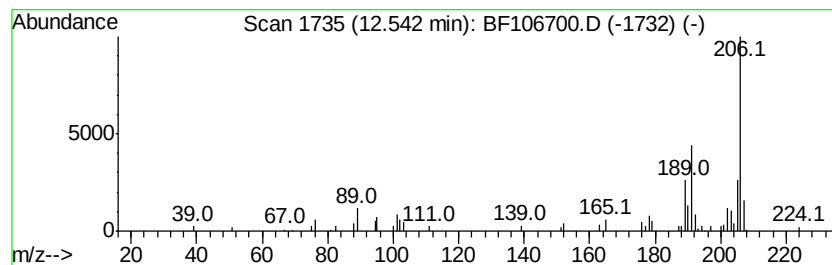
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 9,10-Dimethylantracene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.54	2.05 ng	68393	Phenanthrene-d10		11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene			206	C16H14	000781-43-1	93
2	di-p-Tolylacetylene			206	C16H14	002789-88-0	93
3	Phenanthrene, 2,5-dimethyl-			206	C16H14	003674-66-6	93
4	Anthracene, 1,4-dimethyl-			206	C16H14	000781-92-0	89
5	Phenanthrene, 2,3-dimethyl-			206	C16H14	003674-65-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

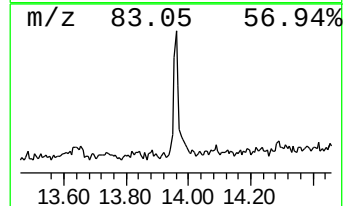
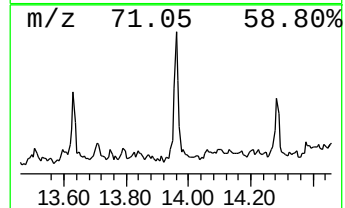
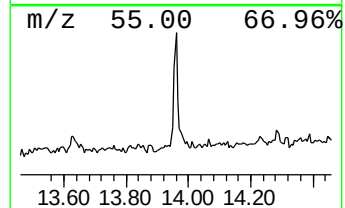
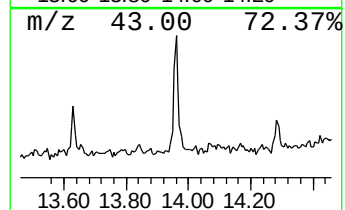
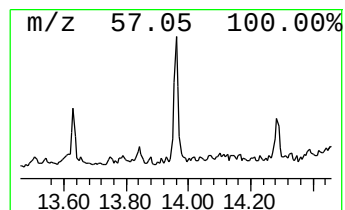
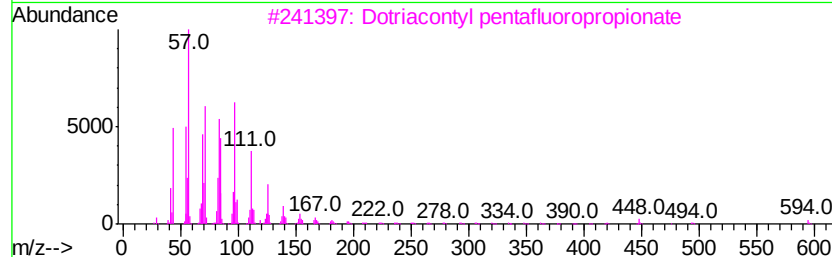
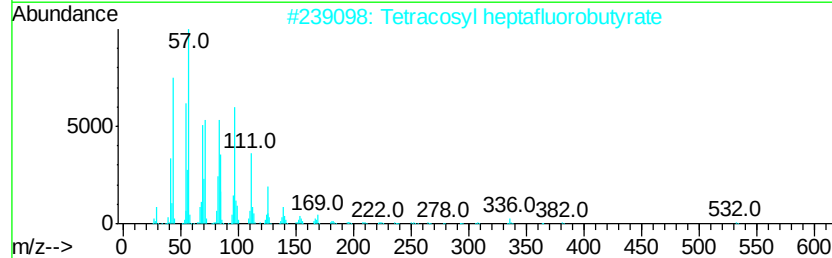
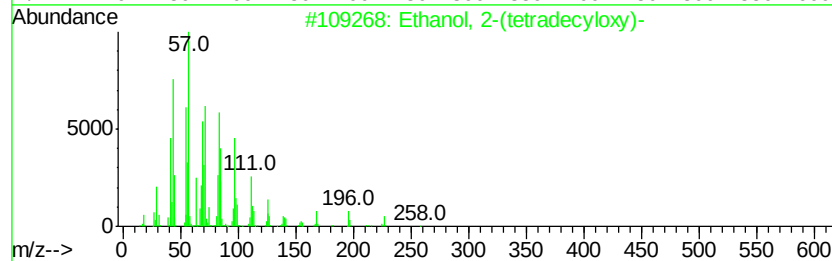
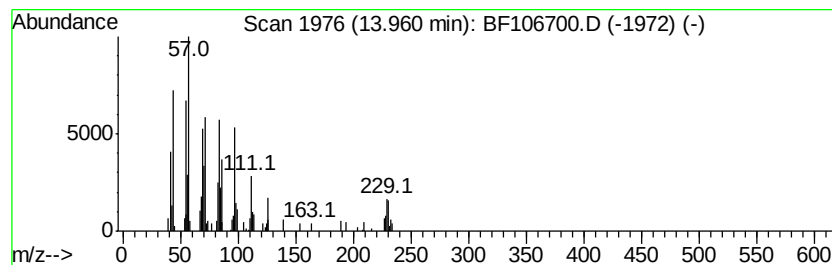
Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.96	2.32 ng	68821	Chrysene-d12		14.15	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
2		Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	81
3		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	74
4		Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	74
5		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062218\
Data File : BF106700.D
Acq On : 22 Jun 2018 11:18
Operator : JU/SJ
Sample : J3603-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26KV-TP-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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
2-Pentanone, 4-hy...	5.20	7.9	ng	161151	1	6.96	407605	20.0
unknown6.74	6.74	74.1	ng	1510230	1	6.96	407605	20.0
Hexadecane	10.88	2.3	ng	76743	4	11.49	665694	20.0
9,10-Dimethylanthe...	12.54	2.0	ng	68393	4	11.49	665694	20.0
Ethanol, 2-(tetra...	13.96	2.3	ng	68821	5	14.15	594224	20.0