

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
 Data File : BF091432.D
 Acq On : 5 Dec 2016 21:50
 Operator : UM/SJ
 Sample : H5838-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-17

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.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	2.355	3	6	17	rVB3	34009	97559	1.67%	0.252%
2	4.516	192	195	198	rBV	348905	392090	6.72%	1.013%
3	5.019	237	239	242	rBV	355109	524551	8.99%	1.355%
4	5.441	274	276	279	rBV	1987068	1989821	34.12%	5.140%
5	6.470	364	366	368	rBV	1485896	1251386	21.46%	3.233%
6	6.619	376	379	381	rBV	3684476	4135350	70.91%	10.682%
7	6.847	396	399	401	rBV	1093151	926920	15.89%	2.394%
8	7.007	410	413	415	rVB	3724067	3579366	61.38%	9.246%
9	7.407	445	448	451	rBV	2171262	3213212	55.10%	8.300%
10	8.127	509	511	513	rVB	973716	1159836	19.89%	2.996%
11	8.802	568	570	572	rVB	68562	67756	1.16%	0.175%
12	8.905	577	579	584	rVV2	40683	88525	1.52%	0.229%
13	9.213	602	606	608	rBV	5586171	5646292	96.82%	14.585%
14	9.888	662	665	667	rVB	1695597	1515803	25.99%	3.916%
15	10.402	705	710	711	rBV4	46193	62312	1.07%	0.161%
16	10.676	731	734	736	rBV	3878985	4068803	69.77%	10.510%
17	10.768	741	742	747	rBV2	26685	76109	1.31%	0.197%
18	11.259	781	785	787	rBV	78986	96344	1.65%	0.249%
19	11.362	792	794	797	rVV	1088582	1407669	24.14%	3.636%
20	11.899	839	841	843	rVB	129597	108307	1.86%	0.280%
21	12.688	908	910	916	rBV	117135	152276	2.61%	0.393%
22	12.802	916	920	922	rVB	71528	87527	1.50%	0.226%
23	12.962	931	934	936	rBV	4797399	5831586	100.00%	15.064%
24	14.002	1022	1025	1027	rBV	1123147	1298143	22.26%	3.353%
25	15.420	1147	1149	1152	rBV	654635	934929	16.03%	2.415%

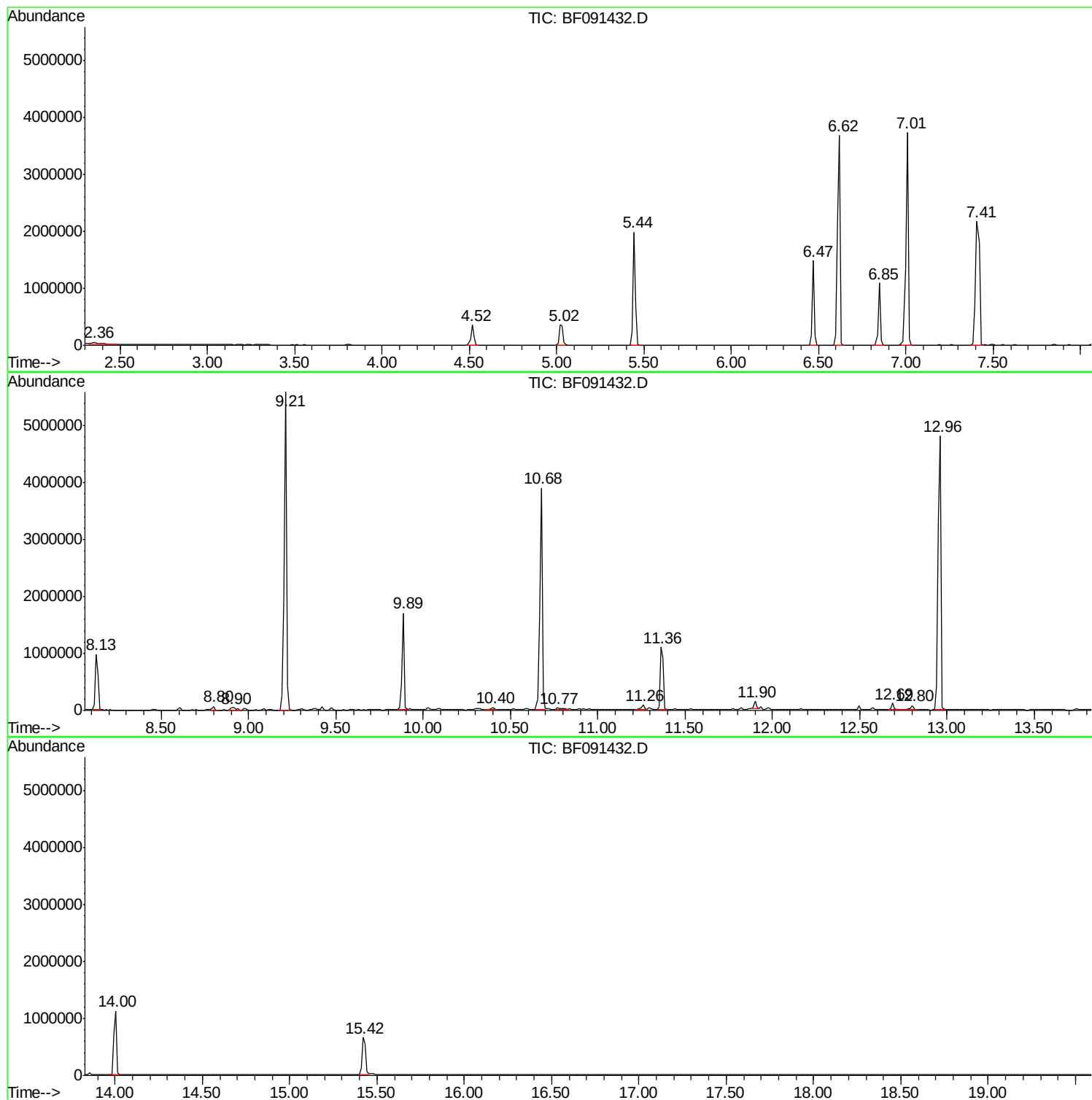
Sum of corrected areas: 38712472

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
Data File : BF091432.D
Acq On : 5 Dec 2016 21:50
Operator : UM/SJ
Sample : H5838-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-17

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120516\
Data File : BF091432.D
Acq On : 5 Dec 2016 21:50
Operator : UM/SJ
Sample : H5838-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-17

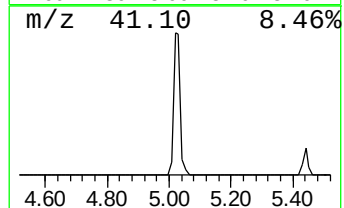
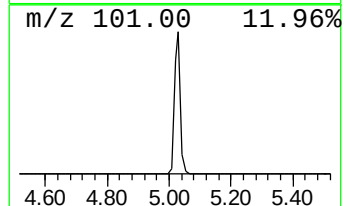
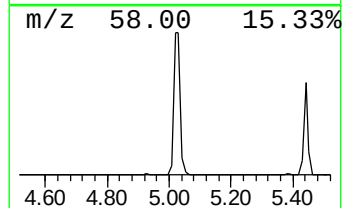
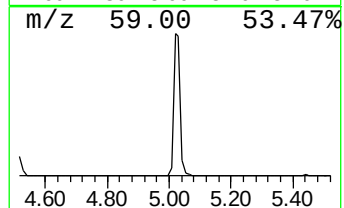
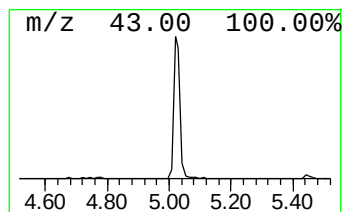
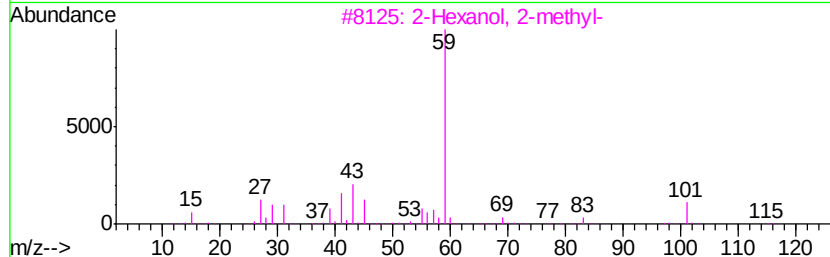
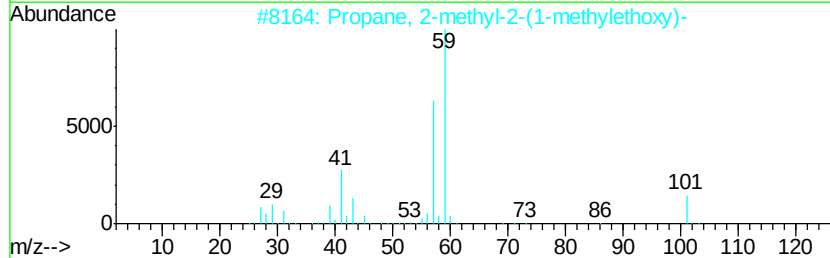
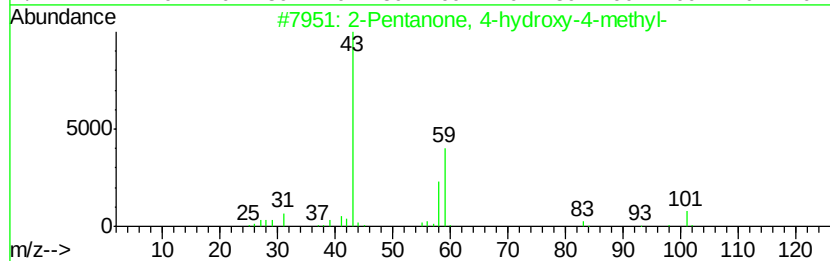
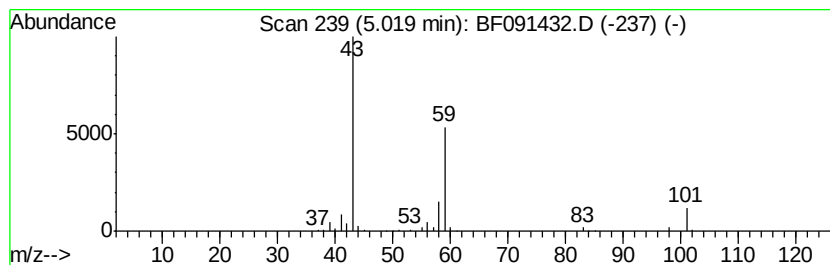
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	11.32 ng	524551	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	33
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
Data File : BF091432.D
Acq On : 5 Dec 2016 21:50
Operator : UM/SJ
Sample : H5838-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-17

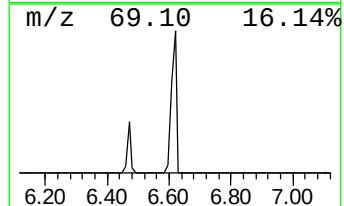
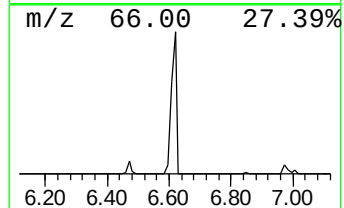
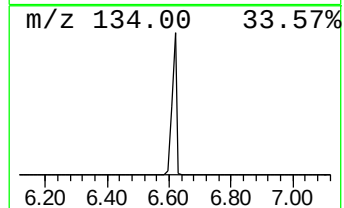
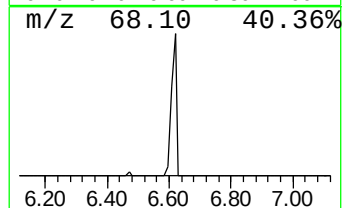
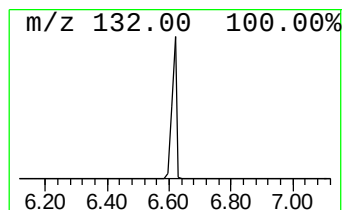
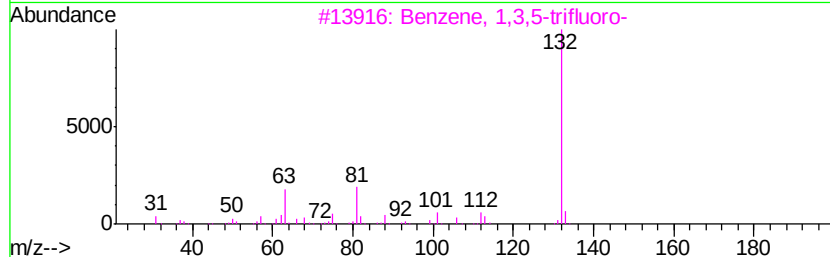
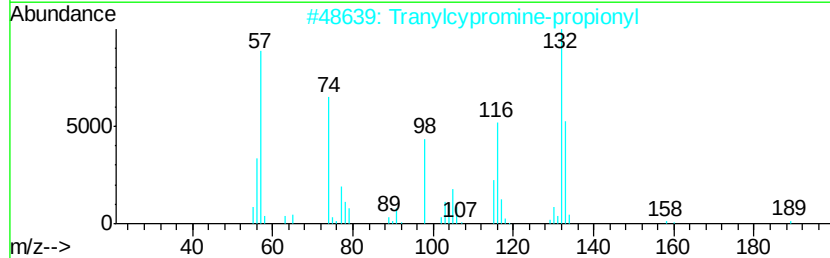
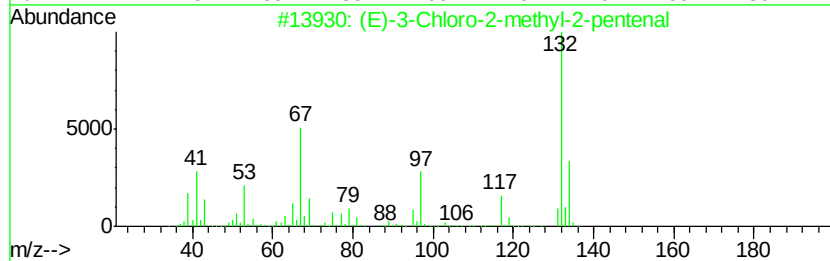
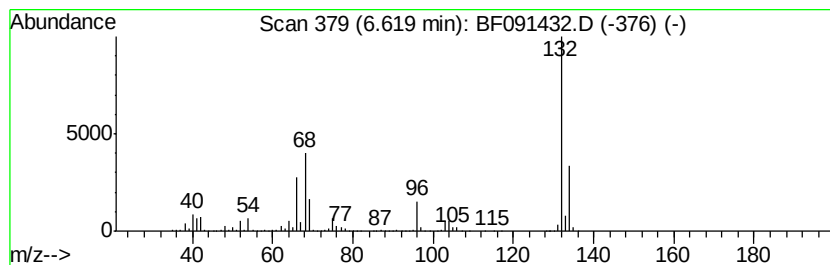
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	89.23 ng	4135350	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
Data File : BF091432.D
Acq On : 5 Dec 2016 21:50
Operator : UM/SJ
Sample : H5838-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

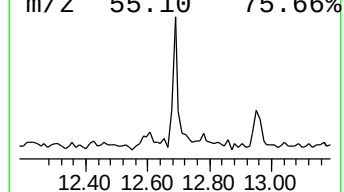
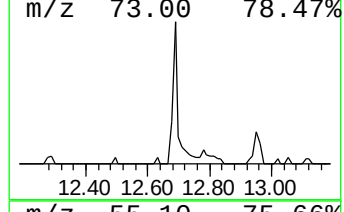
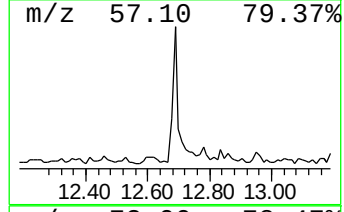
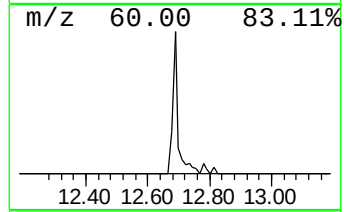
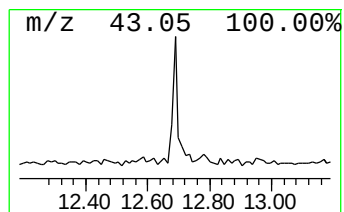
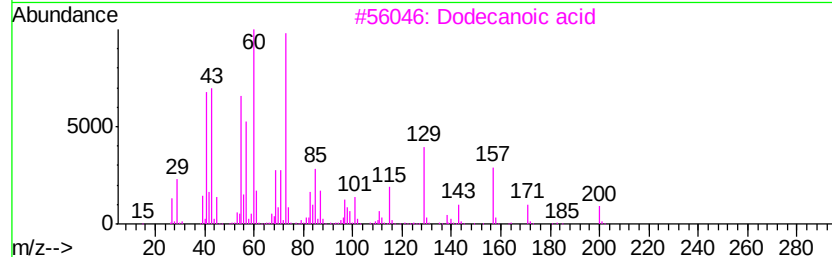
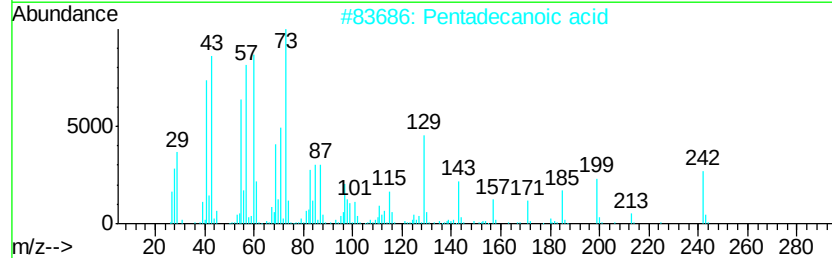
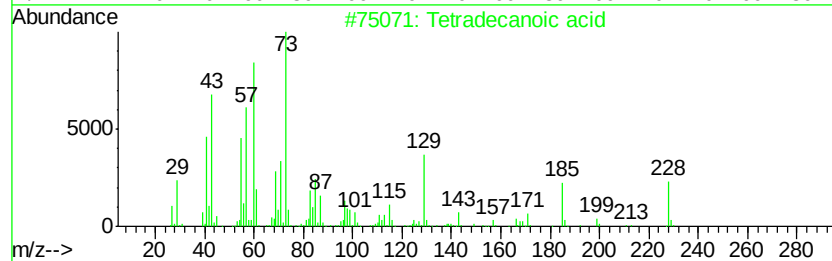
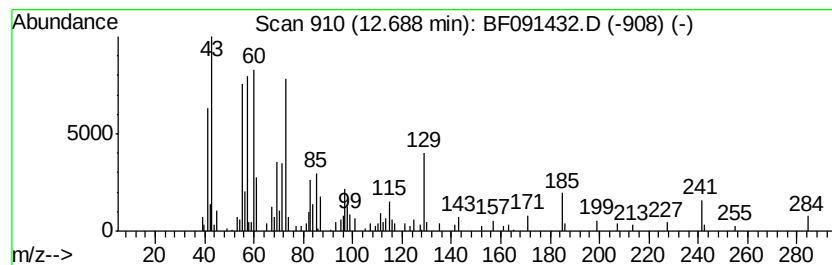
Instrument :
BNA_F
ClientSampleId :
MW-17

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetradecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.69	2.35 ng	152276	Chrysene-d12		14.00
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
3	Dodecanoic acid	200	C12H24O2	000143-07-7	64
4	Undecanoic acid	186	C11H22O2	000112-37-8	50
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
Data File : BF091432.D
Acq On : 5 Dec 2016 21:50
Operator : UM/SJ
Sample : H5838-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-17

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	11.3	ng	524551	1	6.85	926920	20.0
unknown6.62	6.62	89.2	ng	4135350	1	6.85	926920	20.0
Tetradecanoic acid	12.69	2.4	ng	152276	5	14.00	1298140	20.0