

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062918\
 Data File : BF106998.D
 Acq On : 30 Jun 2018 2:51
 Operator : JU/SJ
 Sample : J3718-06
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-16

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.195	482	486	495	rBV	209429	246828	11.70%	1.606%
2	5.595	550	554	568	rBV	1383314	1402349	66.47%	9.123%
3	6.595	720	724	736	rBV	1134518	1177047	55.79%	7.657%
4	6.731	743	747	750	rBV	1796391	1673304	79.32%	10.885%
5	6.948	781	784	788	rBV	478509	405628	19.23%	2.639%
6	7.107	807	811	815	rBV	1553986	1445867	68.54%	9.406%
7	7.519	876	881	884	rBV	1142340	1161291	55.05%	7.554%
8	8.236	999	1003	1017	rVB	491864	502852	23.84%	3.271%
9	9.307	1180	1185	1188	rBV	2194270	2109657	100.00%	13.724%
10	9.701	1245	1252	1256	rBV	47932	47993	2.27%	0.312%
11	9.989	1298	1301	1308	rVB	560890	544386	25.80%	3.541%
12	10.401	1368	1371	1374	rVB	34347	25677	1.22%	0.167%
13	10.789	1432	1437	1440	rBV	1372579	1266075	60.01%	8.236%
14	10.871	1448	1451	1457	rVB	27602	27944	1.32%	0.182%
15	11.483	1552	1555	1559	rBV	558427	504878	23.93%	3.284%
16	13.071	1820	1825	1828	rBV	2004968	1870320	88.66%	12.167%
17	13.942	1970	1973	1978	rVB	44438	43712	2.07%	0.284%
18	14.136	2002	2006	2017	rBV	551161	513695	24.35%	3.342%
19	15.671	2262	2267	2277	rVB	296973	402715	19.09%	2.620%

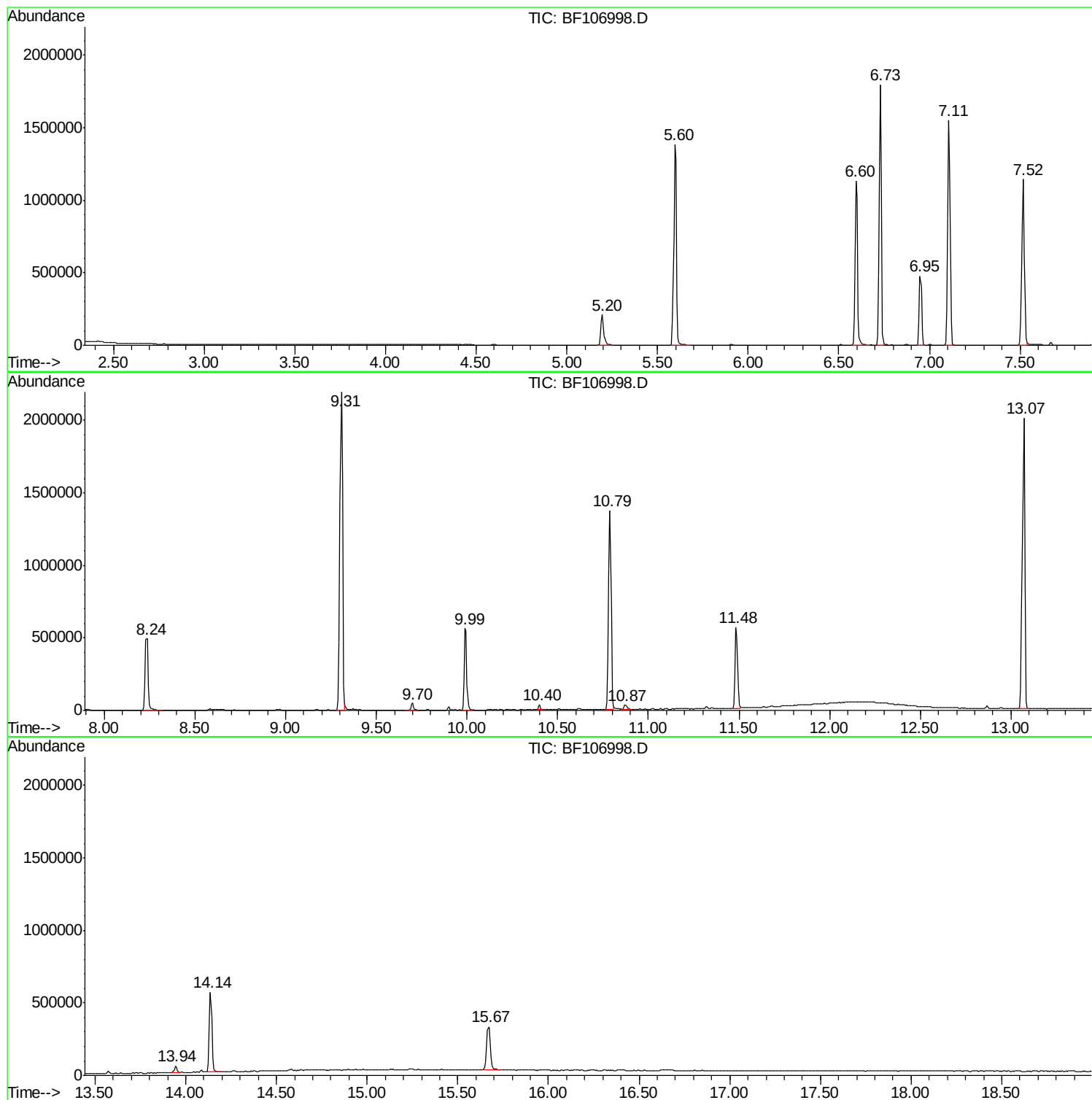
Sum of corrected areas: 15372218

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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



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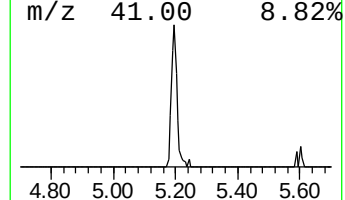
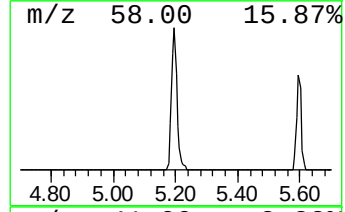
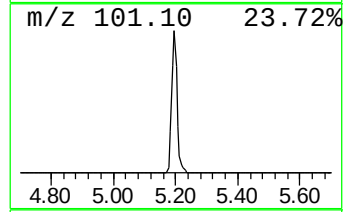
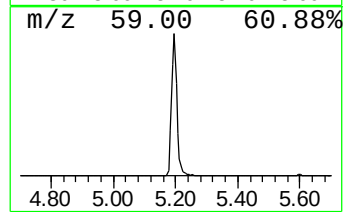
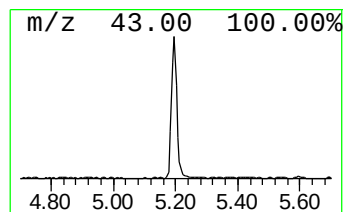
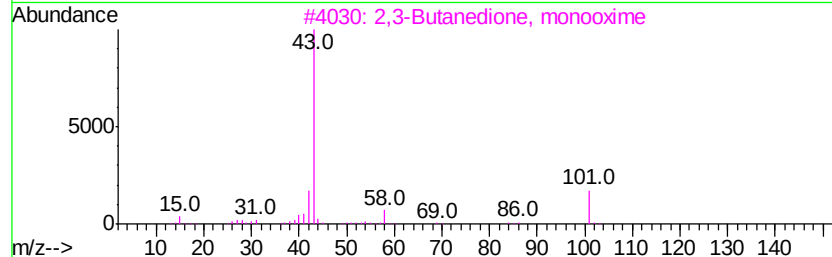
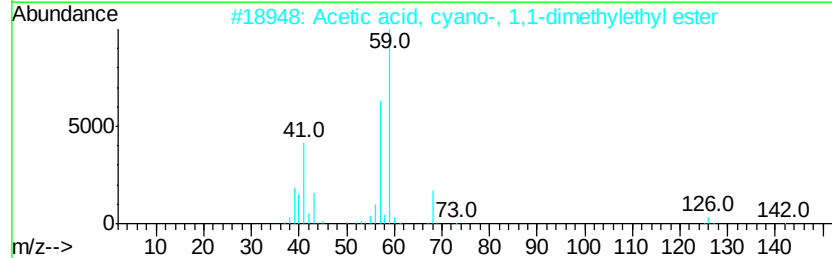
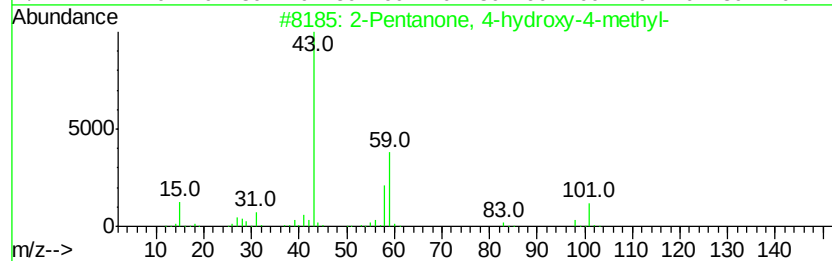
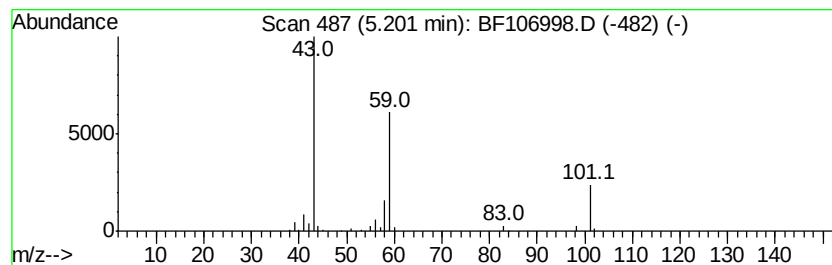
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	12.17 ng	246828	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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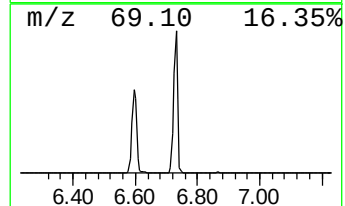
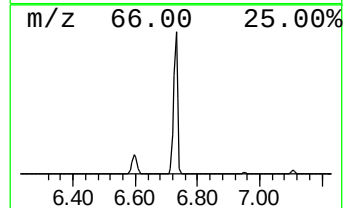
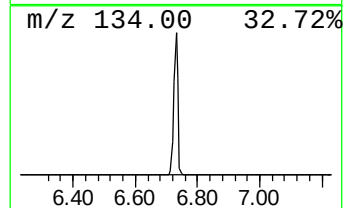
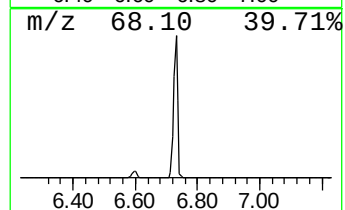
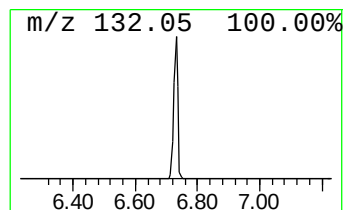
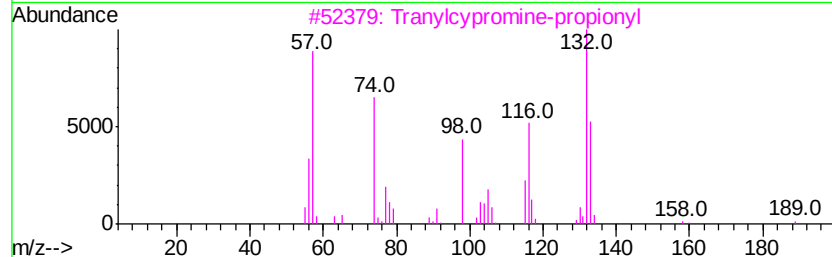
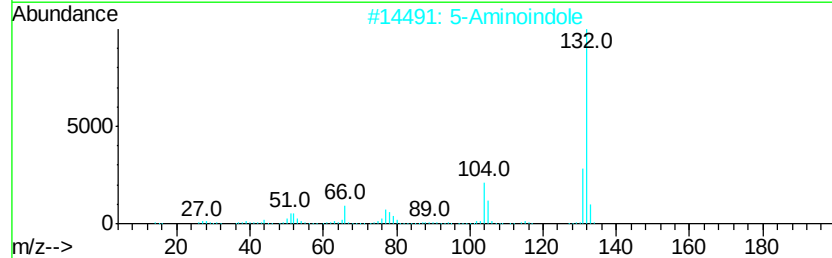
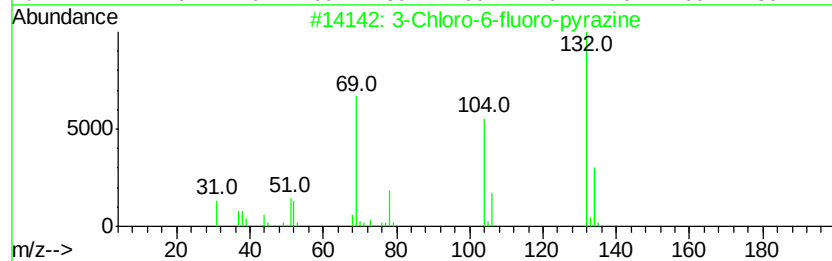
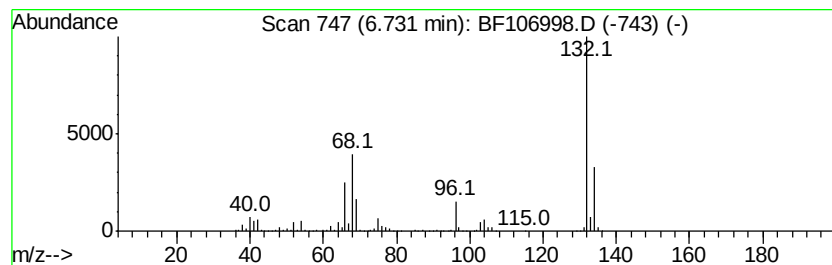
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	82.50 ng	1673300	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2	5		Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	12.2	ng	246828	1	6.95	405628	20.0
unknown6.73	6.73	82.5	ng	1673300	1	6.95	405628	20.0