

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112817\
 Data File : BF100960.D
 Acq On : 29 Nov 2017 3:57
 Operator : SJ/JU
 Sample : I6603-02
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112217-TIDO-F-D-S

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:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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	5.075	504	508	516	rBV	88099	95170	5.29%	0.969%
2	5.475	572	576	584	rBV	500454	432396	24.04%	4.402%
3	6.481	744	747	755	rVB	320022	275615	15.32%	2.806%
4	6.628	768	772	775	rBV	1019445	843817	46.92%	8.591%
5	6.857	808	811	815	rVB	451099	355232	19.75%	3.617%
6	7.016	834	838	841	rBV	1535762	1277792	71.05%	13.009%
7	7.422	902	907	910	rBV	938518	966876	53.76%	9.844%
8	8.139	1025	1029	1033	rBV	546181	430074	23.91%	4.379%
9	9.216	1207	1212	1215	rBV	2072822	1798484	100.00%	18.310%
10	9.898	1324	1328	1338	rVB	454517	425905	23.68%	4.336%
11	10.680	1458	1461	1473	rBV	560814	544457	30.27%	5.543%
12	11.380	1577	1580	1584	rBV	443554	365100	20.30%	3.717%
13	12.968	1845	1850	1853	rBV	1524747	1372554	76.32%	13.974%
14	13.862	1997	2002	2007	rBV2	15008	22229	1.24%	0.226%
15	14.021	2026	2029	2034	rBV	395284	338583	18.83%	3.447%
16	15.492	2275	2279	2286	rBV	204754	277966	15.46%	2.830%

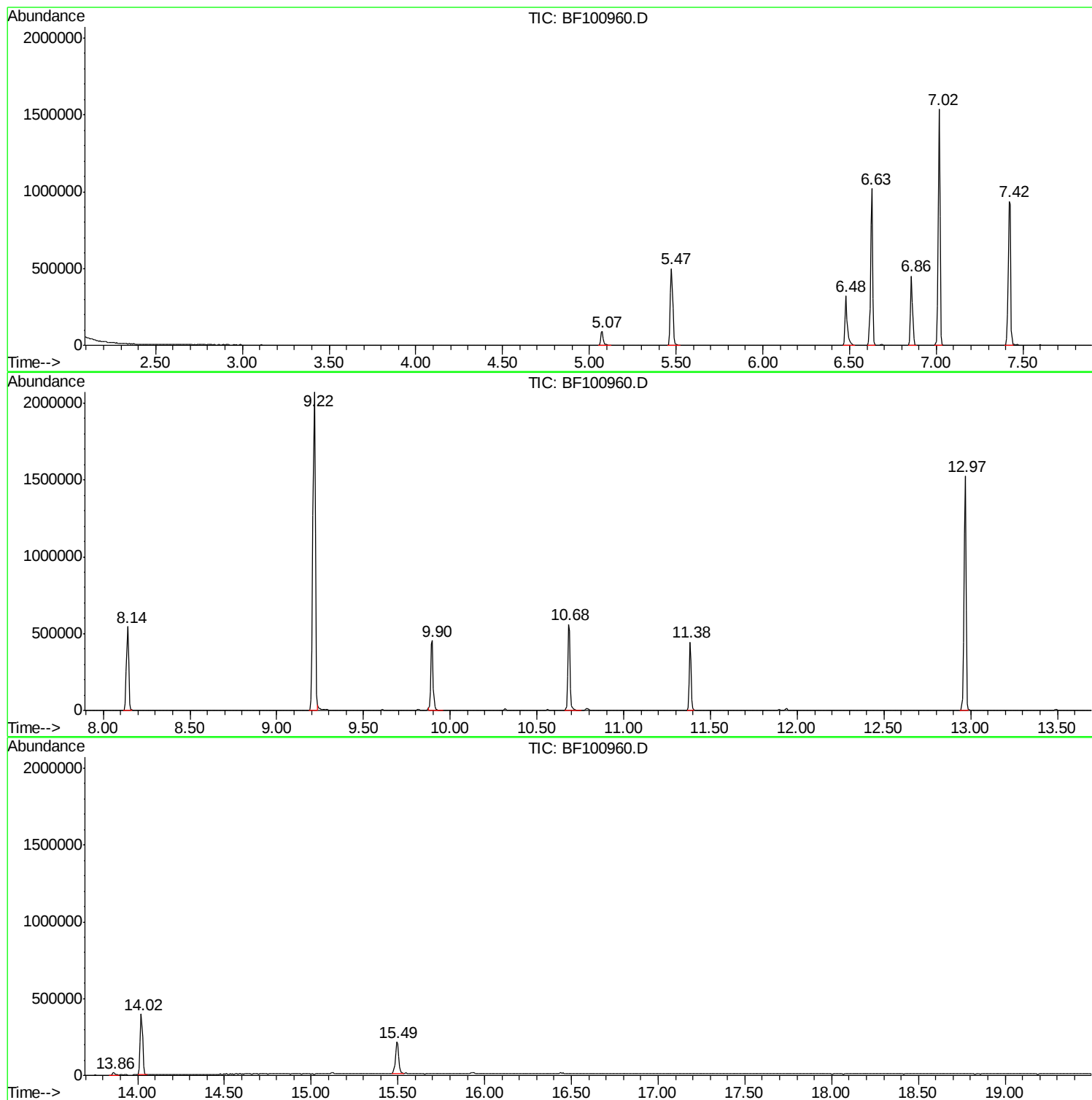
Sum of corrected areas: 9822250

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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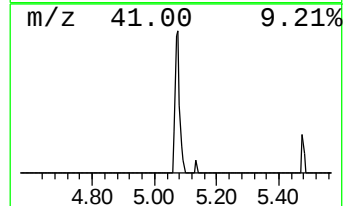
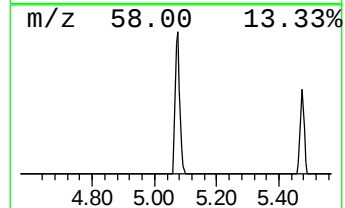
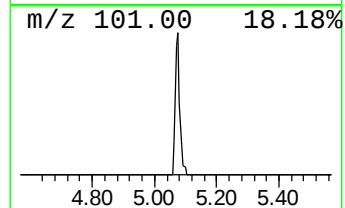
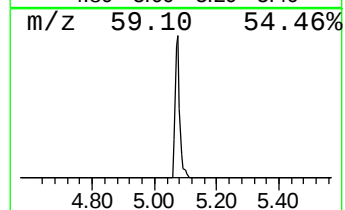
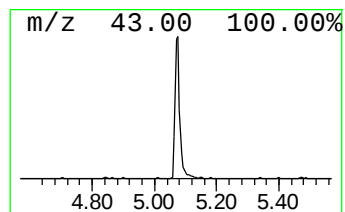
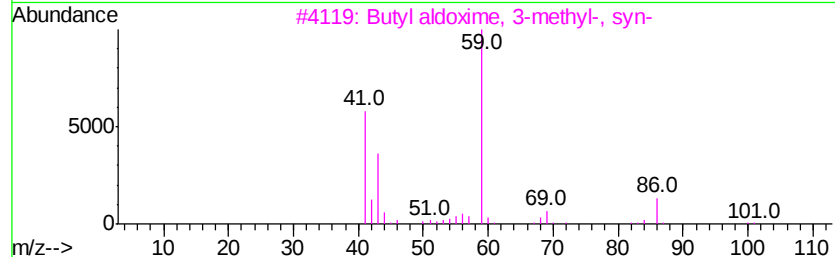
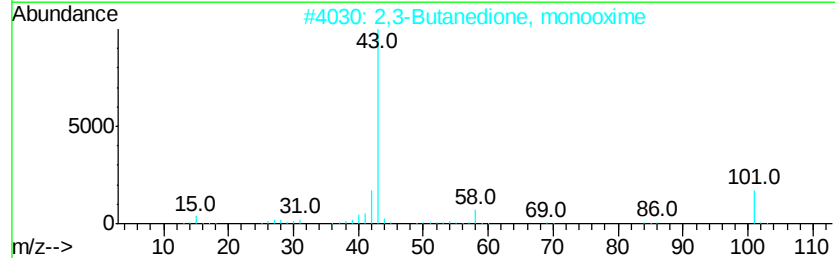
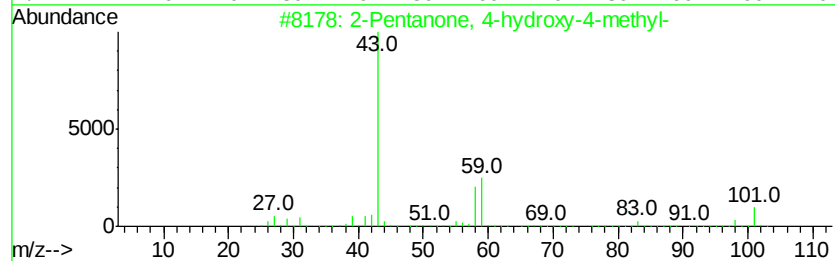
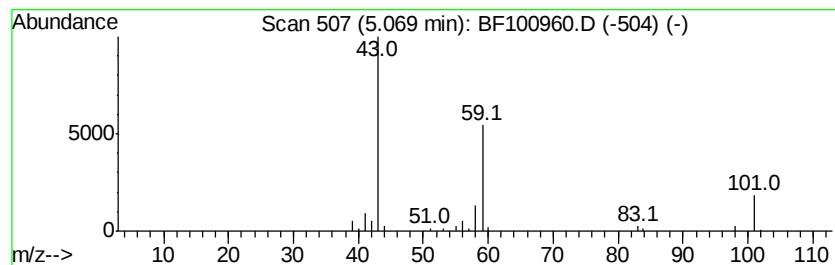
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.36 ng	95170	1,4-Dichlorobenzene-d4	6.86

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	38
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	9



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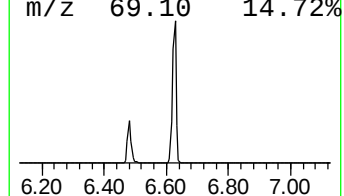
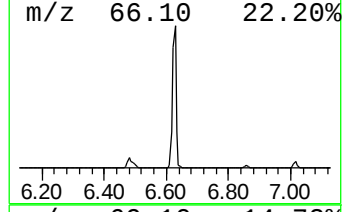
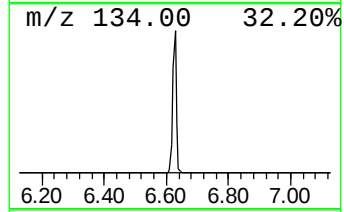
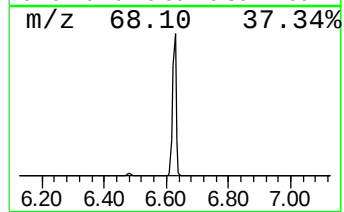
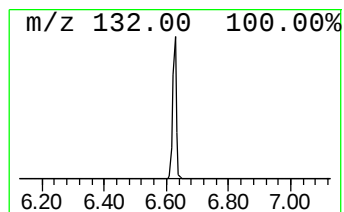
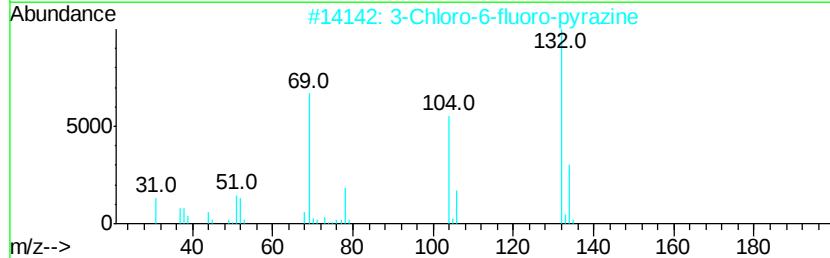
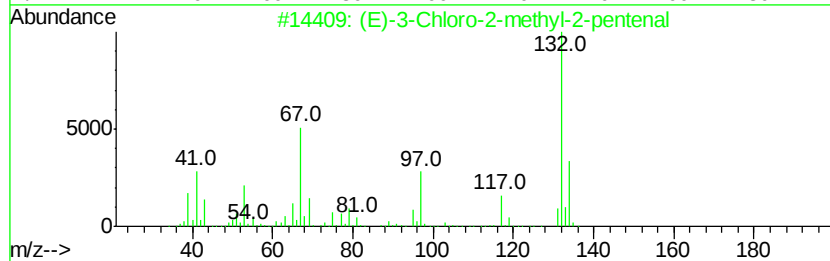
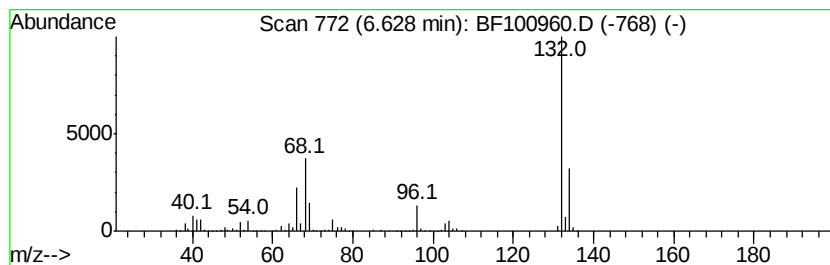
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	47.51 ng	843817	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	5.4	ng	95170	1	6.86	355232	20.0
unknown6.63	6.63	47.5	ng	843817	1	6.86	355232	20.0