

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122017\
 Data File : BF101590.D
 Acq On : 21 Dec 2017 5:39
 Operator : SJ/JU
 Sample : I6971-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121517-TIDO-F-D-B

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-BF121417.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	5.010	494	497	509	rBV	88758	136970	3.69%	0.571%
2	5.416	563	566	585	rBV	1205143	1608276	43.32%	6.704%
3	6.434	736	739	758	rBV	739183	1114415	30.02%	4.646%
4	6.569	758	762	768	rBV	2262388	2342761	63.11%	9.766%
5	6.792	797	800	811	rBV	926447	884773	23.83%	3.688%
6	6.951	823	827	842	rBV	2995554	2970622	80.02%	12.383%
7	7.363	893	897	922	rBV	2226833	2592124	69.82%	10.806%
8	8.075	1015	1018	1035	rBV	1103402	1183409	31.88%	4.933%
9	9.151	1196	1201	1204	rBV	3960003	3712347	100.00%	15.475%
10	9.833	1313	1317	1321	rBV	1006098	946765	25.50%	3.947%
11	9.863	1321	1322	1330	rVB	56254	44378	1.20%	0.185%
12	10.628	1448	1452	1466	rBV	1204182	1292416	34.81%	5.388%
13	11.322	1566	1570	1573	rBV	917948	838709	22.59%	3.496%
14	12.904	1835	1839	1842	rBV	2851171	2599568	70.02%	10.837%
15	13.792	1986	1990	1999	rBV4	30890	60865	1.64%	0.254%
16	13.968	2016	2020	2027	rBV	850413	823885	22.19%	3.434%
17	15.433	2265	2269	2278	rVB	650607	836305	22.53%	3.486%

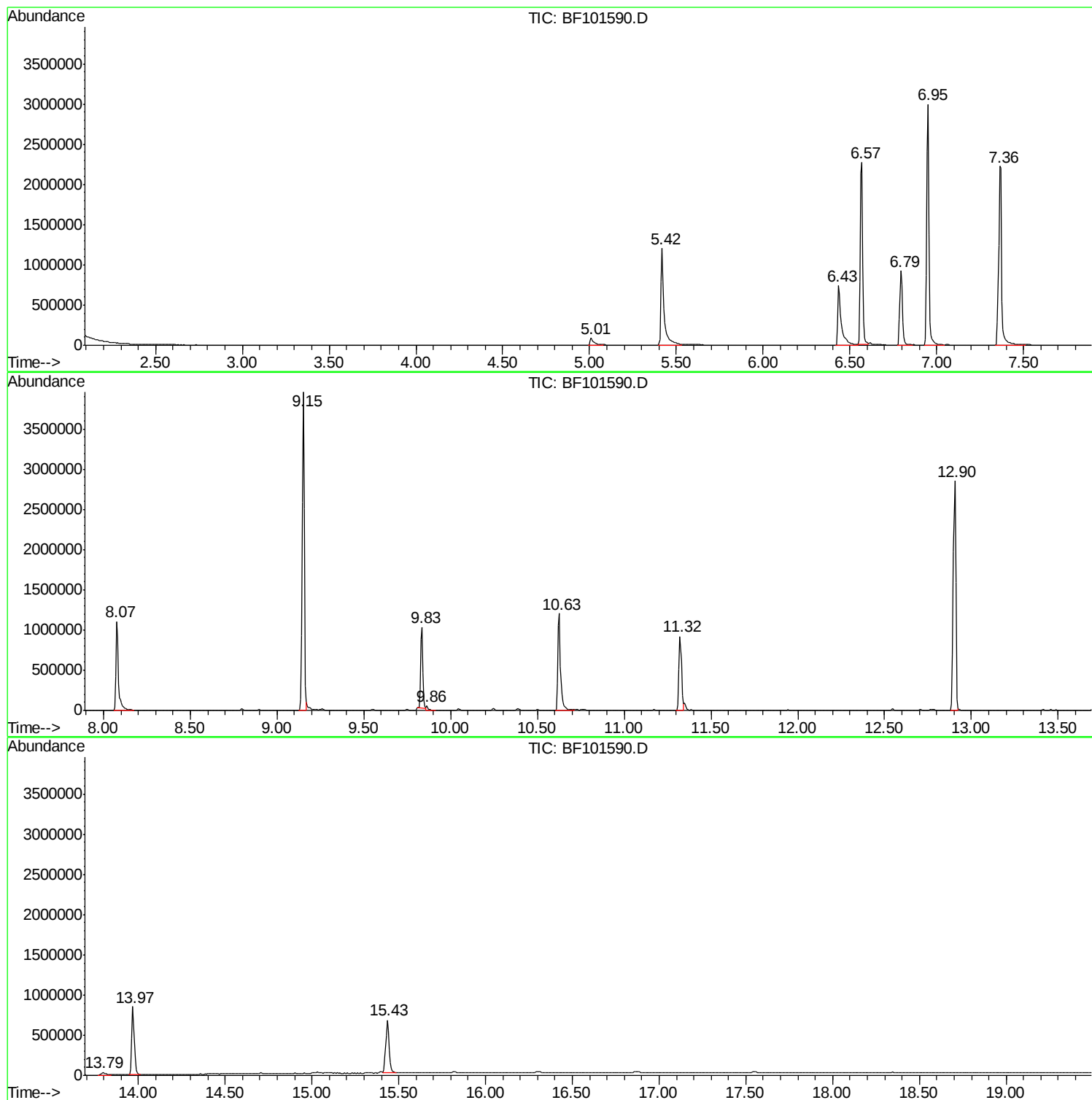
Sum of corrected areas: 23988588

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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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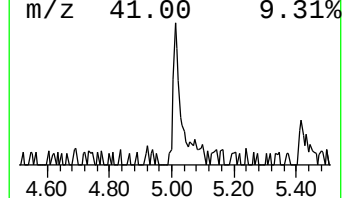
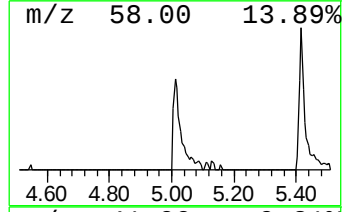
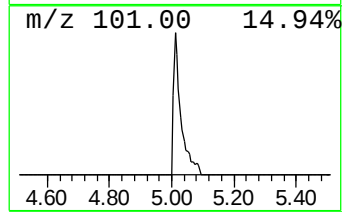
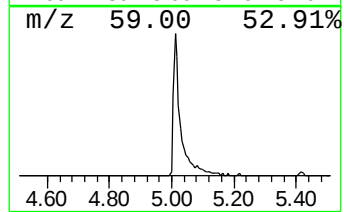
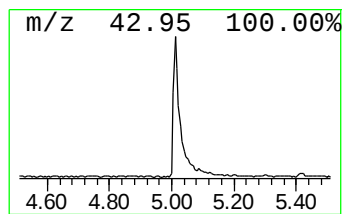
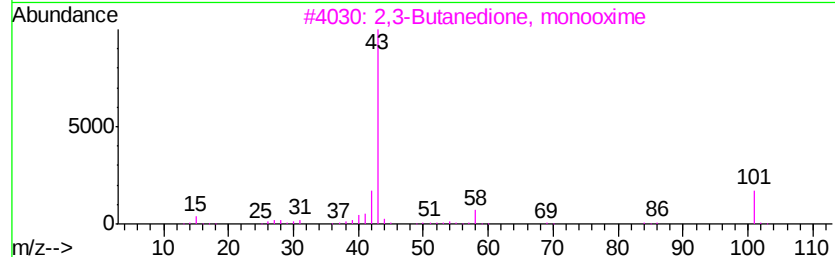
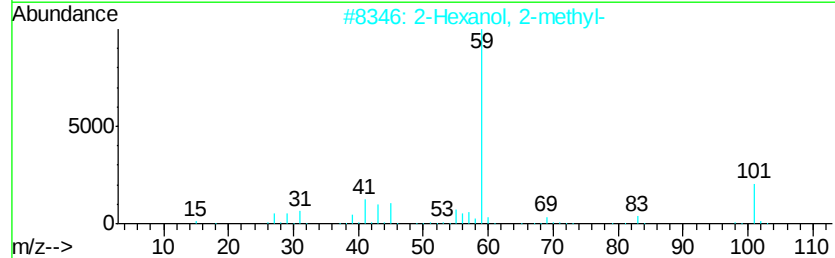
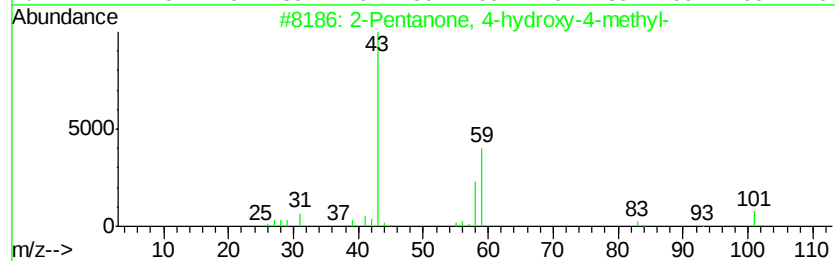
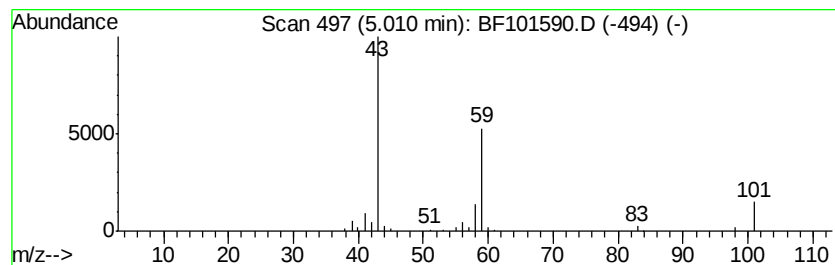
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	3.10 ng	136970	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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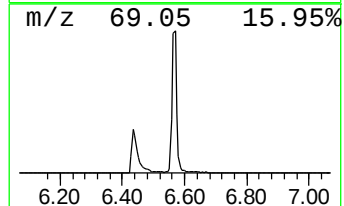
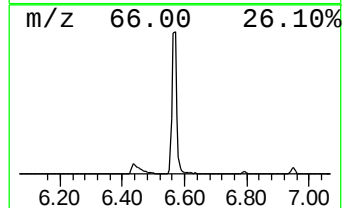
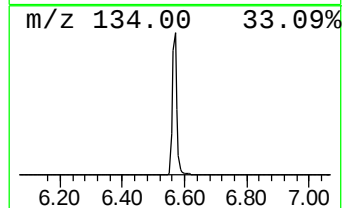
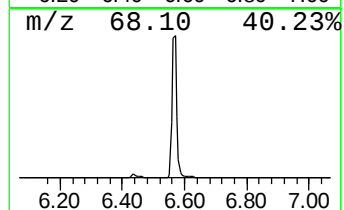
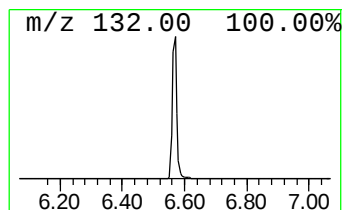
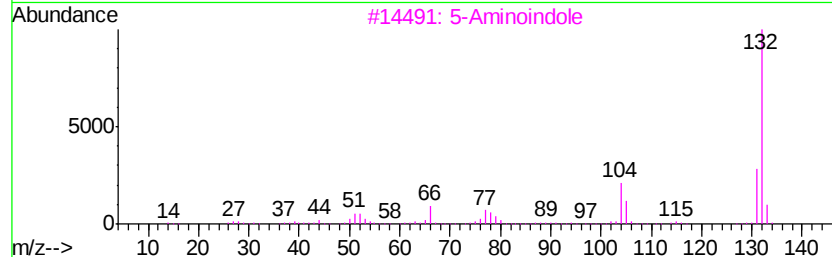
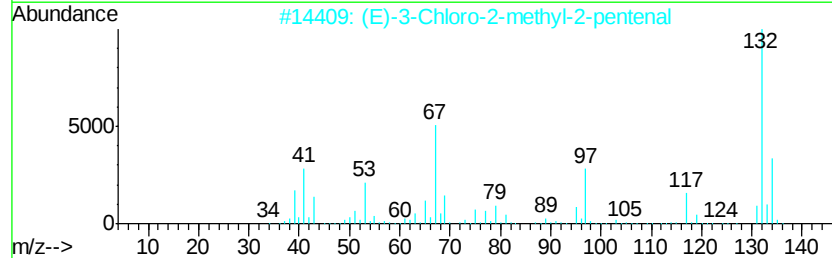
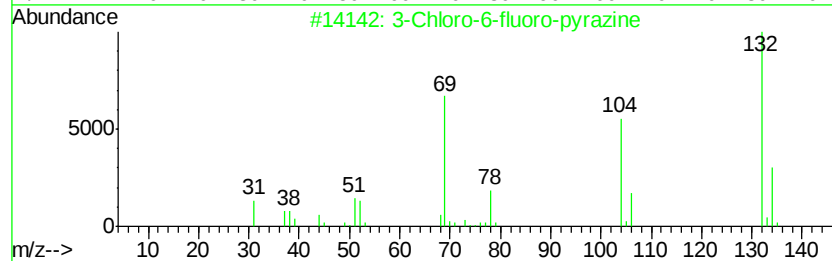
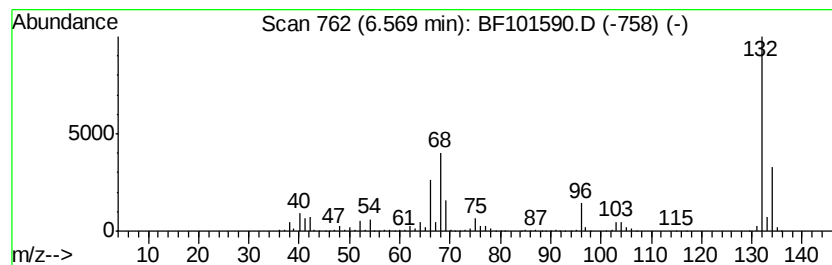
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Peak Number 2 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	52.96 ng	2342760	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	3.1	ng	136970	1	6.79	884773	20.0
unknown6.57	6.57	53.0	ng	2342760	1	6.79	884773	20.0