

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101508.D
 Acq On : 19 Dec 2017 5:11
 Operator : SJ/JU
 Sample : I6909-10
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121217-FOPI-E-U-M

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	5.022	496	499	508	rBV	68722	93082	3.02%	0.470%
2	5.434	566	569	588	rBV	855572	1125428	36.46%	5.683%
3	6.457	740	743	761	rBV	558887	816382	26.45%	4.122%
4	6.581	761	764	776	rVV	1775343	1881346	60.96%	9.500%
5	6.810	799	803	811	rBV	797568	709149	22.98%	3.581%
6	6.963	825	829	846	rBV	2617021	2591484	83.97%	13.086%
7	7.381	895	900	917	rBV	1972451	2025706	65.63%	10.229%
8	8.092	1017	1021	1036	rBV	859728	828826	26.85%	4.185%
9	9.169	1199	1204	1207	rBV	3175774	3086333	100.00%	15.585%
10	9.563	1268	1271	1276	rVB	84365	71802	2.33%	0.363%
11	9.845	1316	1319	1330	rVB	722051	708565	22.96%	3.578%
12	10.563	1437	1441	1446	rBV	34524	34376	1.11%	0.174%
13	10.645	1451	1455	1462	rBV	950501	998939	32.37%	5.044%
14	11.339	1570	1573	1582	rBV	673383	651608	21.11%	3.290%
15	11.851	1656	1660	1665	rBV	47197	63939	2.07%	0.323%
16	12.563	1776	1781	1787	rBV8	23545	51209	1.66%	0.259%
17	12.921	1838	1842	1846	rBV	2585494	2491966	80.74%	12.583%
18	13.804	1989	1992	2001	rBV	60551	73385	2.38%	0.371%
19	13.992	2020	2024	2029	rBV	802383	788814	25.56%	3.983%
20	15.462	2270	2274	2283	rVB	527670	711422	23.05%	3.592%

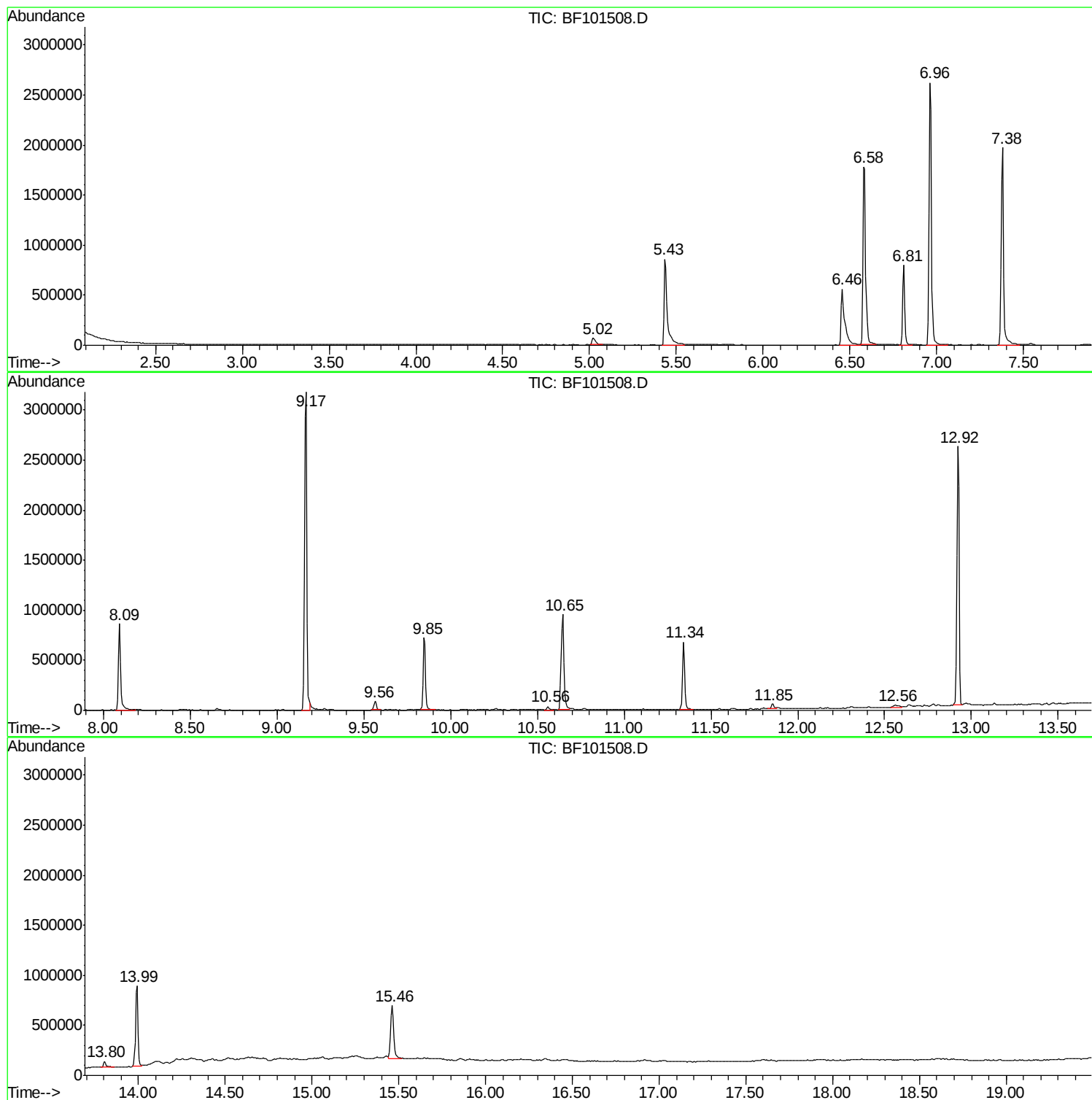
Sum of corrected areas: 19803761

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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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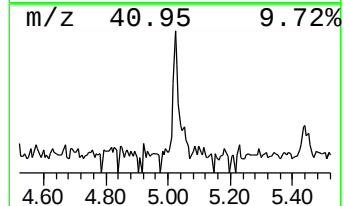
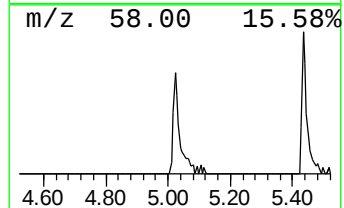
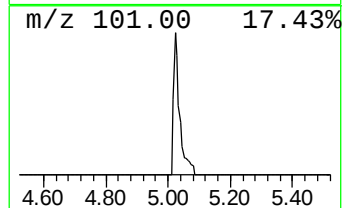
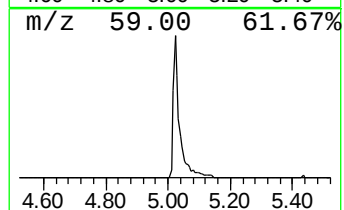
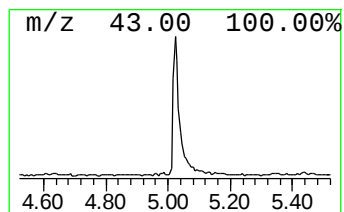
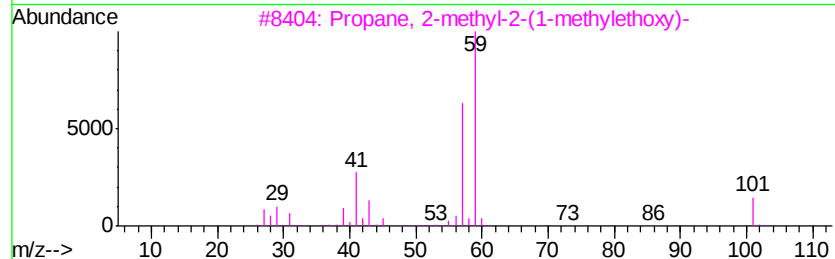
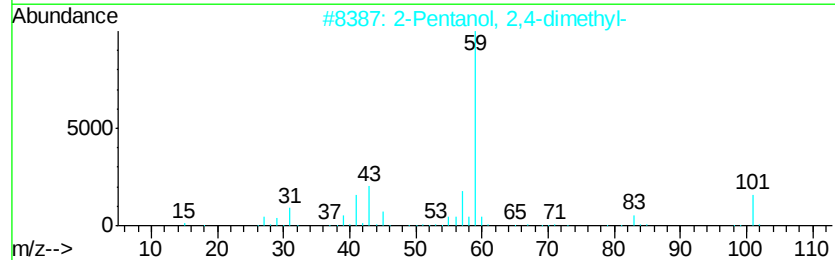
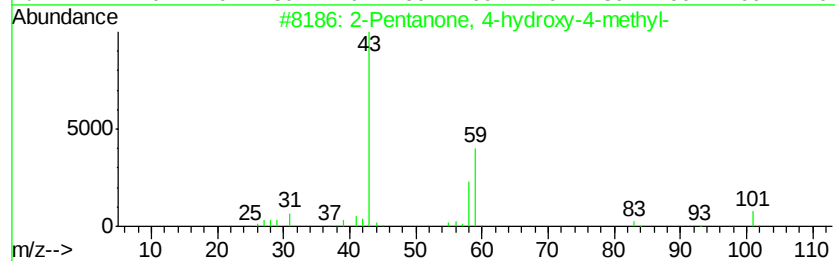
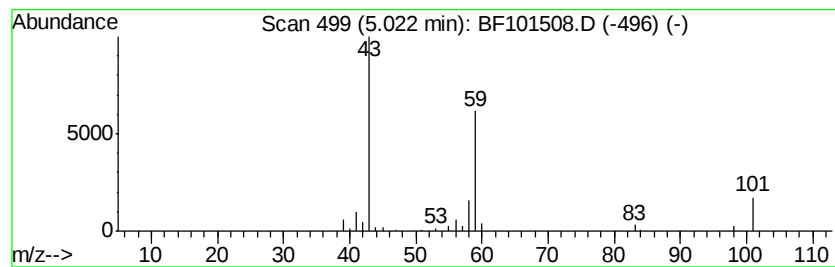
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	2.63 ng	93082	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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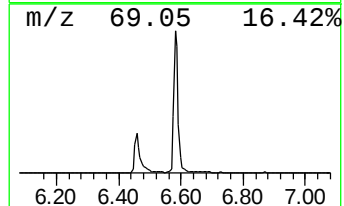
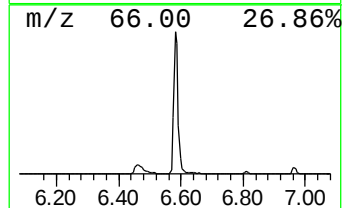
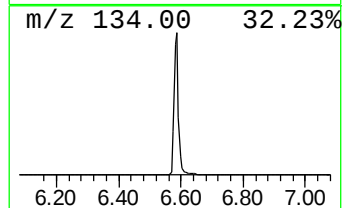
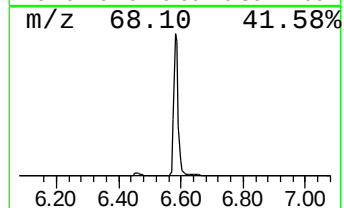
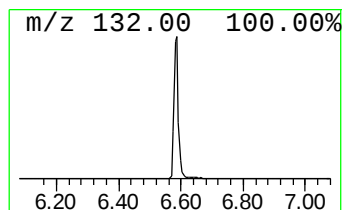
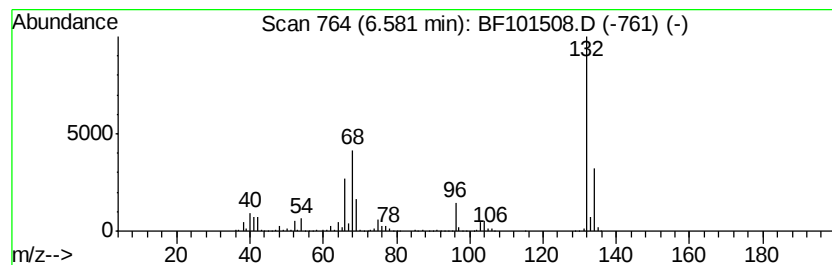
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Peak Number 2 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	53.06 ng	1881350	1,4-Dichlorobenzene-d4	6.81

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			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	2.6	ng	93082	1	6.81	709149	20.0
unknown6.58	6.58	53.1	ng	1881350	1	6.81	709149	20.0