

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122117\
 Data File : BF101634.D
 Acq On : 22 Dec 2017 4:54
 Operator : SJ/JU
 Sample : I7008-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121817-DBRI-F-U-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-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	5.004	493	496	512	rBV	113532	182899	5.69%	0.933%
2	5.416	563	566	594	rBV	668537	1095580	34.06%	5.588%
3	6.440	737	740	758	rBV	417947	729353	22.67%	3.720%
4	6.569	758	762	769	rBV	1705577	1846623	57.41%	9.418%
5	6.793	797	800	810	rBV	736019	750097	23.32%	3.826%
6	6.951	823	827	844	rBV	2542407	2487519	77.33%	12.687%
7	7.363	893	897	913	rBV	1849431	2112385	65.67%	10.773%
8	8.075	1015	1018	1036	rBV	824338	937072	29.13%	4.779%
9	9.151	1197	1201	1211	rBV	3366912	3216683	100.00%	16.405%
10	9.834	1314	1317	1330	rVB	768972	813473	25.29%	4.149%
11	10.628	1448	1452	1465	rBV	865060	994880	30.93%	5.074%
12	11.322	1566	1570	1583	rBV	677250	671349	20.87%	3.424%
13	12.904	1834	1839	1842	rBV	2341017	2183964	67.89%	11.138%
14	13.969	2016	2020	2028	rBV	772082	754399	23.45%	3.848%
15	15.398	2258	2263	2265	rBV	30802	43154	1.34%	0.220%
16	15.433	2265	2269	2286	rVB	414659	617442	19.19%	3.149%
17	15.815	2330	2334	2340	rVB	30666	42757	1.33%	0.218%
18	16.298	2411	2416	2421	rBV	28296	50542	1.57%	0.258%
19	16.868	2508	2513	2520	rVB3	18950	35067	1.09%	0.179%
20	17.539	2621	2627	2635	rVB6	16250	42207	1.31%	0.215%

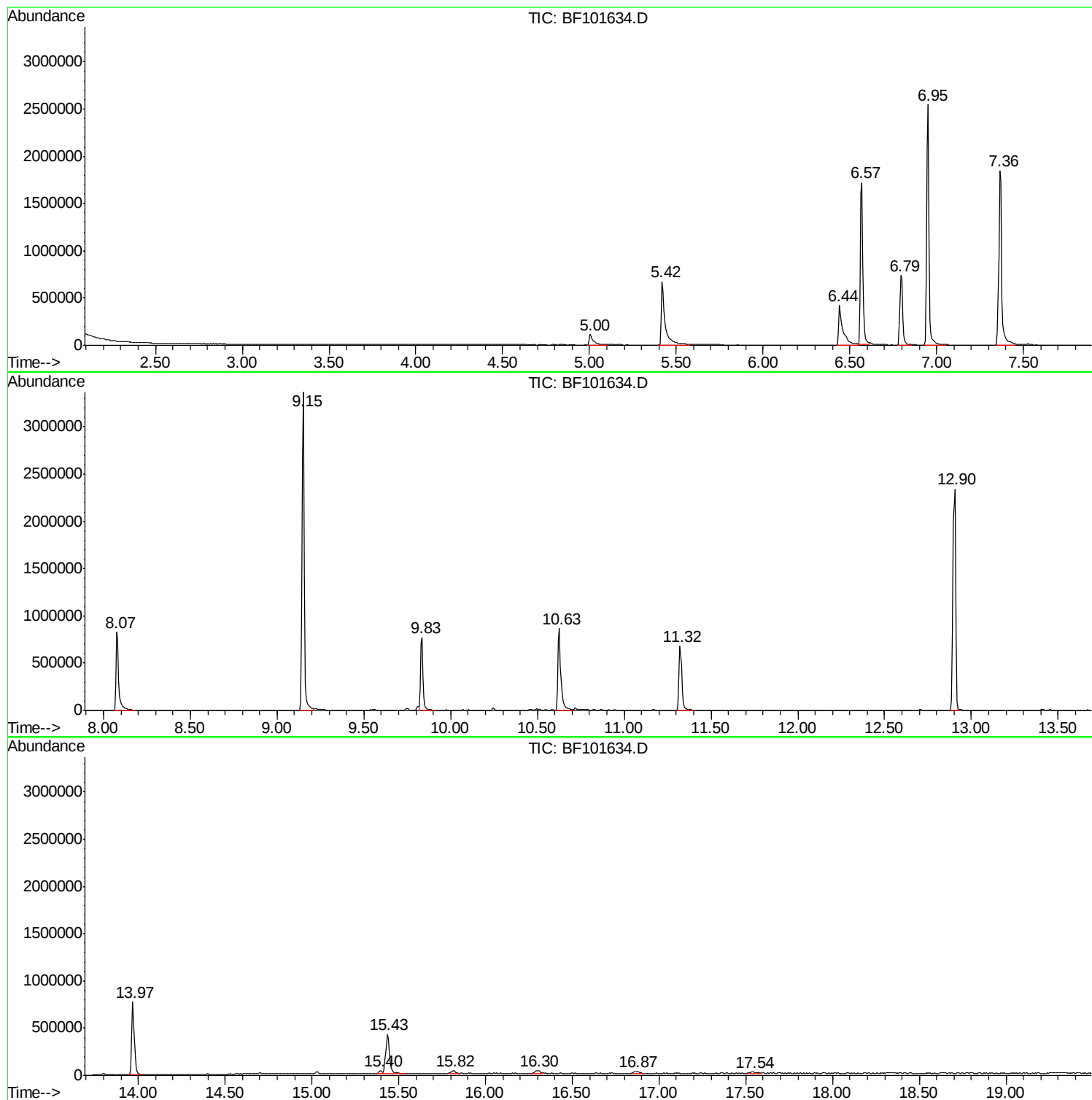
Sum of corrected areas: 19607445

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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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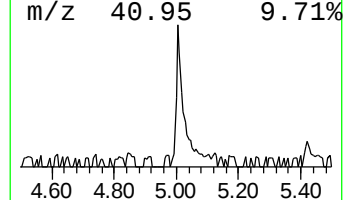
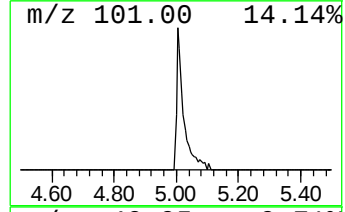
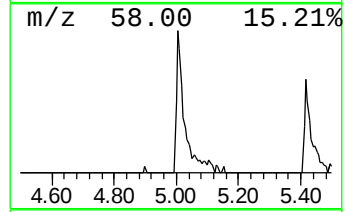
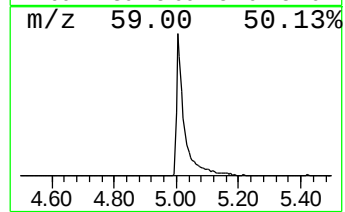
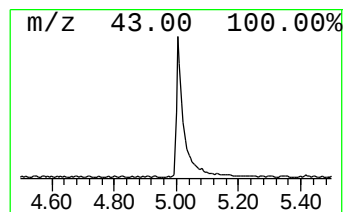
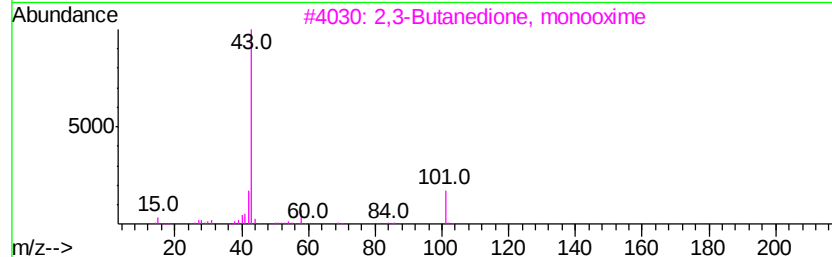
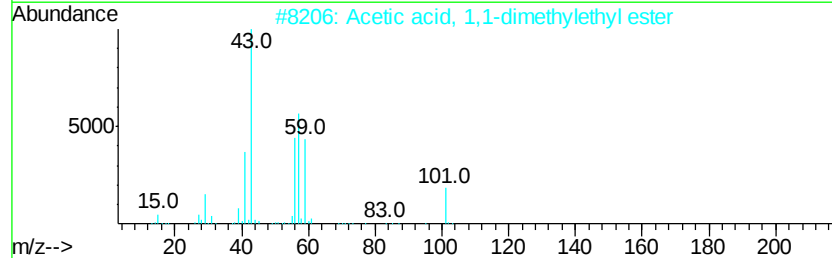
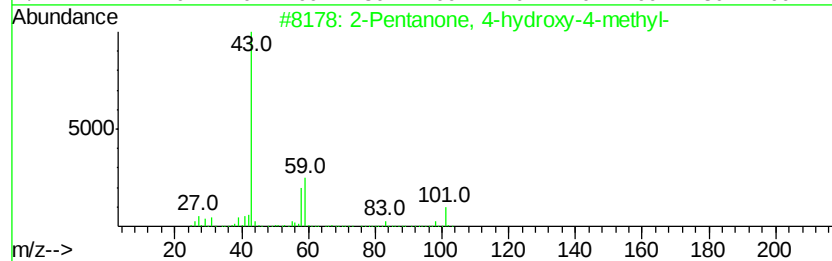
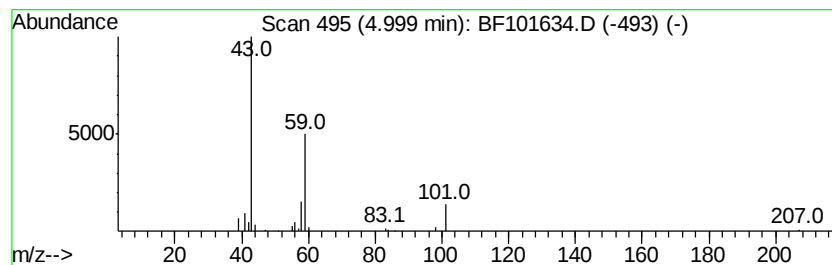
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	4.88 ng	182899	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	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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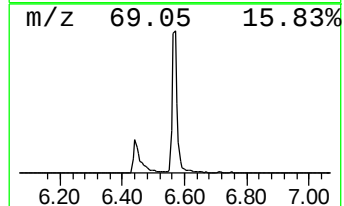
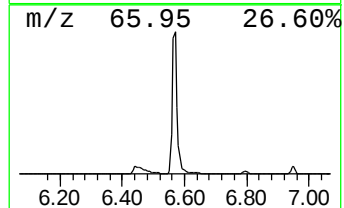
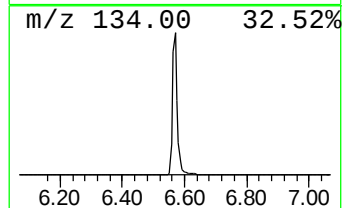
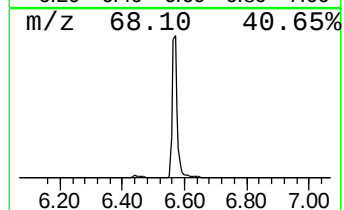
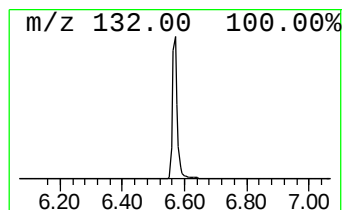
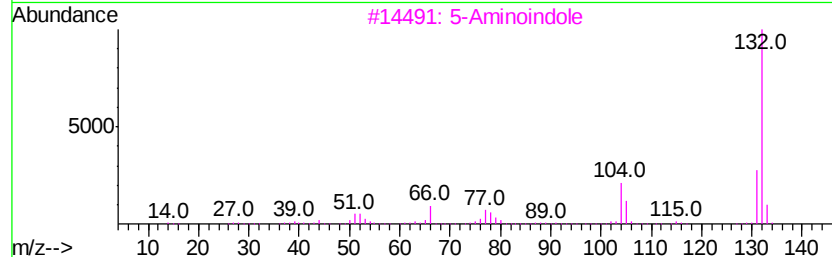
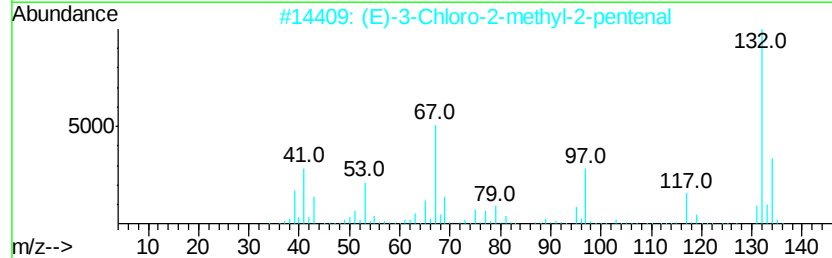
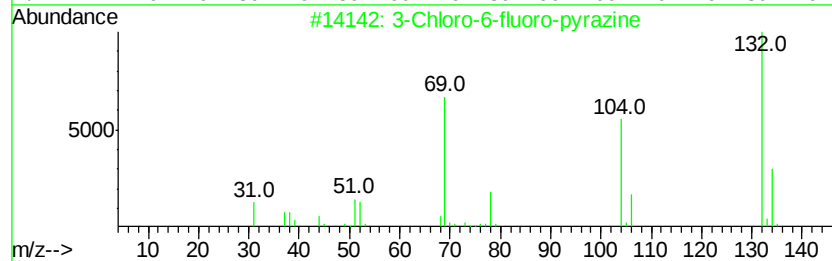
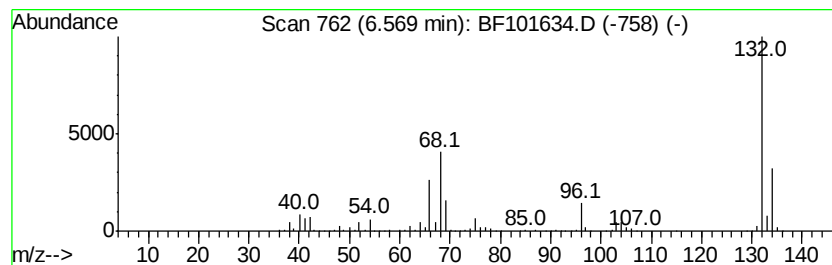
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Peak Number 2 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	49.24 ng	1846620	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	4.9	ng	182899	1	6.79	750097	20.0
unknown6.57	6.57	49.2	ng	1846620	1	6.79	750097	20.0