

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041218\
 Data File : BF104448.D
 Acq On : 12 Apr 2018 14:01
 Operator : JU/SJ
 Sample : PB108173BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB108173BL

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-BF040618.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	5.034	495	501	509	rVB	99423	135264	3.70%	0.488%
2	5.428	563	568	571	rBV	2769517	2866027	78.48%	10.332%
3	6.440	734	740	743	rBV	2625756	2805619	76.83%	10.114%
4	6.581	759	764	767	rBV	2811622	2932913	80.31%	10.573%
5	6.804	799	802	806	rBV	717896	643519	17.62%	2.320%
6	6.963	825	829	832	rBV	2600778	2354223	64.47%	8.487%
7	7.375	893	899	902	rBV	1722951	1870691	51.23%	6.744%
8	8.092	1016	1021	1024	rBV	958922	890453	24.38%	3.210%
9	9.169	1198	1204	1207	rBV	3513402	3518765	96.36%	12.685%
10	9.845	1314	1319	1323	rBV	1075749	1013566	27.75%	3.654%
11	10.639	1448	1454	1457	rBV	2224989	2214509	60.64%	7.983%
12	11.333	1568	1572	1576	rBV	1151430	1054507	28.88%	3.801%
13	12.927	1838	1843	1847	rBV	3312663	3651873	100.00%	13.165%
14	13.980	2018	2022	2026	rVB	1143957	1011853	27.71%	3.648%
15	15.445	2266	2271	2280	rVB	619727	776204	21.25%	2.798%

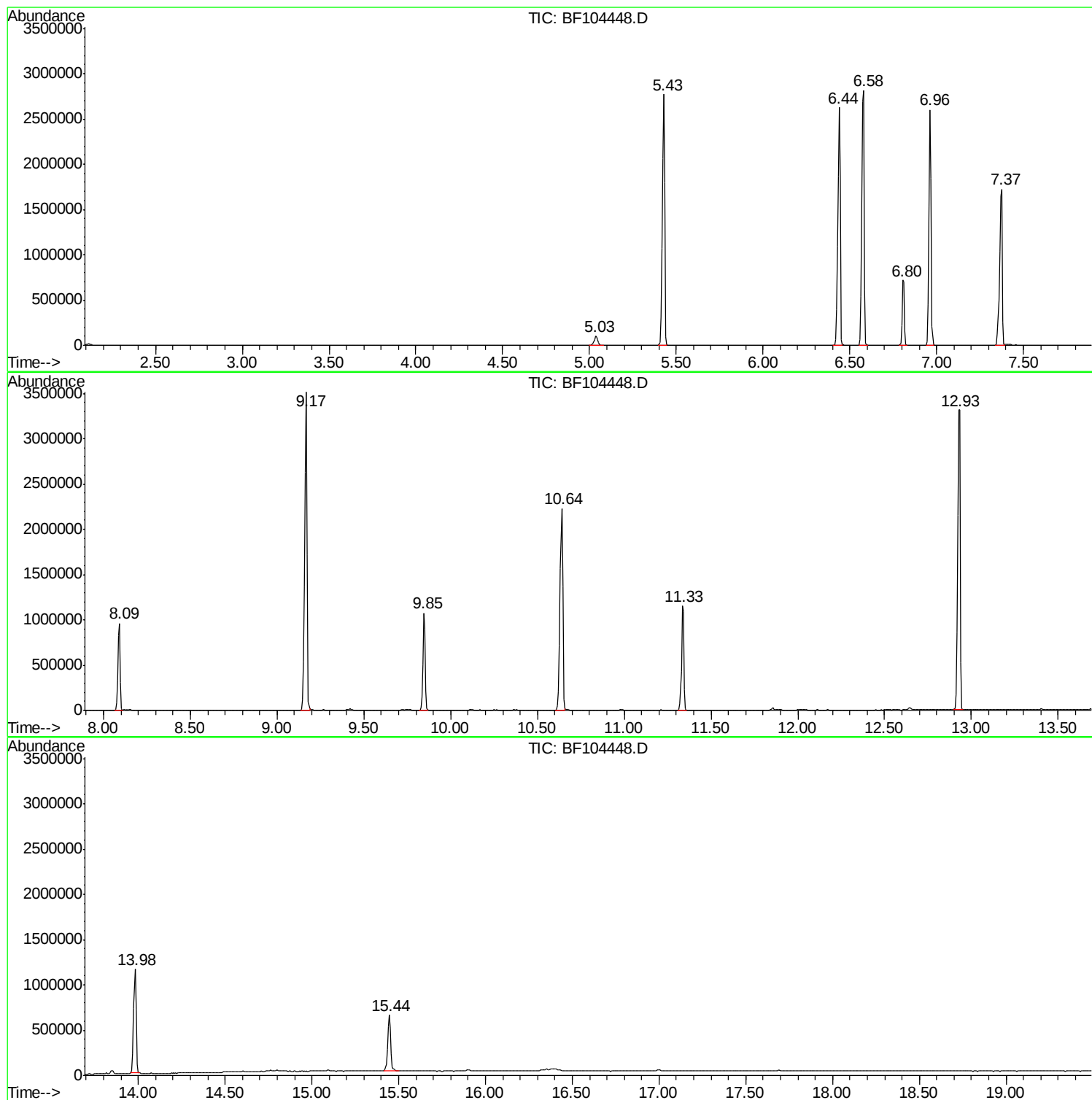
Sum of corrected areas: 27739986

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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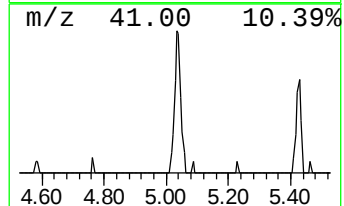
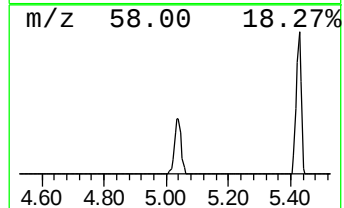
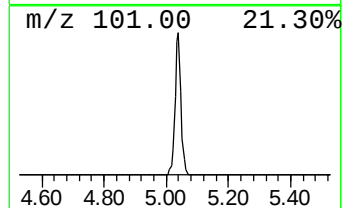
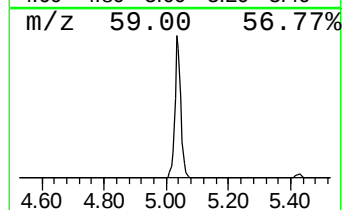
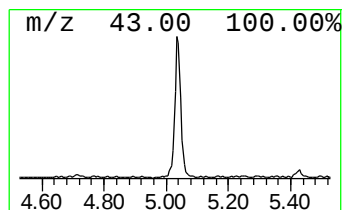
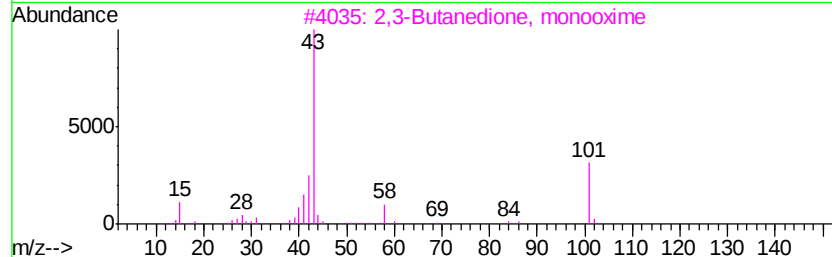
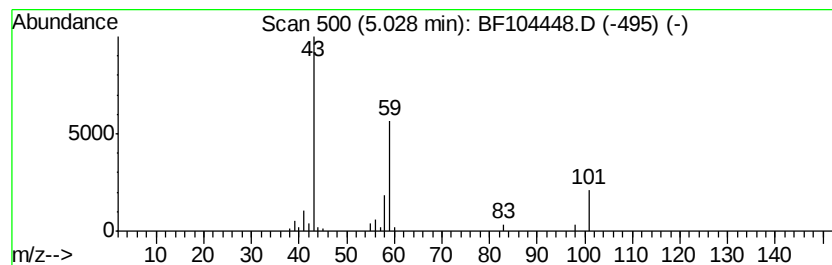
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	4.20 ng	135264	1,4-Dichlorobenzene-d4	6.81

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	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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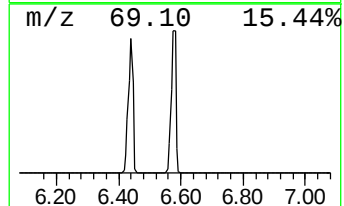
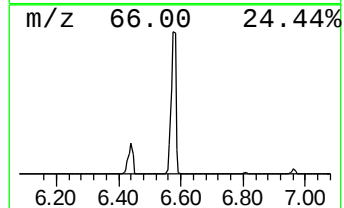
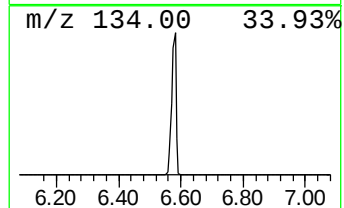
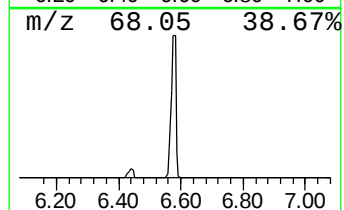
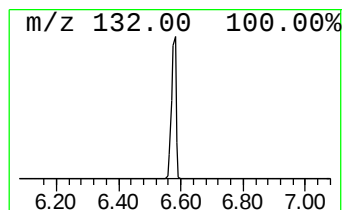
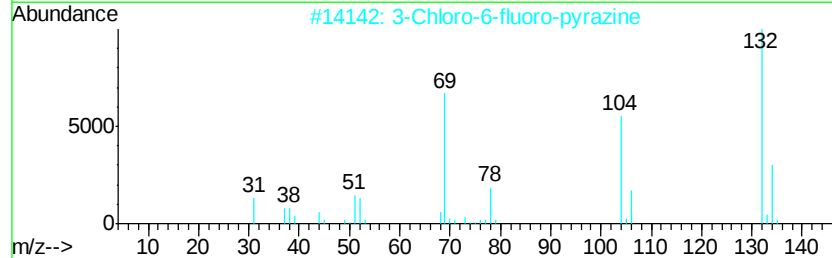
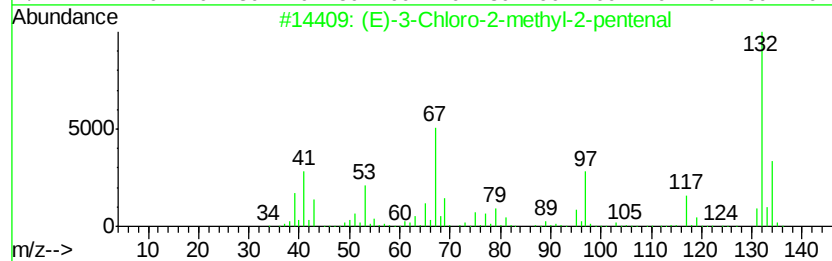
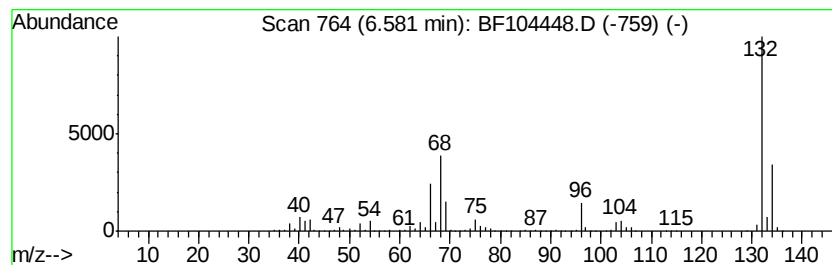
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Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	91.15 ng	2932910	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
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...	5.03	4.2	ng	135264	1	6.81	643519	20.0
unknown6.58	6.58	91.2	ng	2932910	1	6.81	643519	20.0