

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122017\
 Data File : BF101589.D
 Acq On : 21 Dec 2017 5:12
 Operator : SJ/JU
 Sample : PB105190BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB105190BL

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	511	rBV	172550	263033	7.94%	1.087%
2	5.422	563	567	595	rBV	2107295	2569866	77.54%	10.617%
3	6.434	736	739	758	rBV	2147850	2437203	73.54%	10.069%
4	6.569	758	762	775	rVB	2393733	2394840	72.26%	9.894%
5	6.792	797	800	815	rVB	885943	833660	25.15%	3.444%
6	6.951	823	827	845	rBV	2678393	2651581	80.01%	10.955%
7	7.363	893	897	918	rBV	2017943	2197866	66.32%	9.081%
8	8.075	1015	1018	1031	rBV	968905	1019292	30.76%	4.211%
9	9.151	1196	1201	1204	rBV	3435718	3314171	100.00%	13.693%
10	9.833	1312	1317	1332	rBV	913576	942450	28.44%	3.894%
11	10.627	1448	1452	1469	rBV	966405	1053584	31.79%	4.353%
12	11.321	1567	1570	1580	rBV	812238	775756	23.41%	3.205%
13	12.904	1835	1839	1842	rBV	2402312	2160377	65.19%	8.926%
14	13.968	2016	2020	2030	rBV	755924	761662	22.98%	3.147%
15	15.433	2265	2269	2277	rVB	554474	749164	22.60%	3.095%
16	16.304	2412	2417	2423	rBV3	27358	45640	1.38%	0.189%
17	16.868	2509	2513	2520	rVB4	15016	34009	1.03%	0.141%

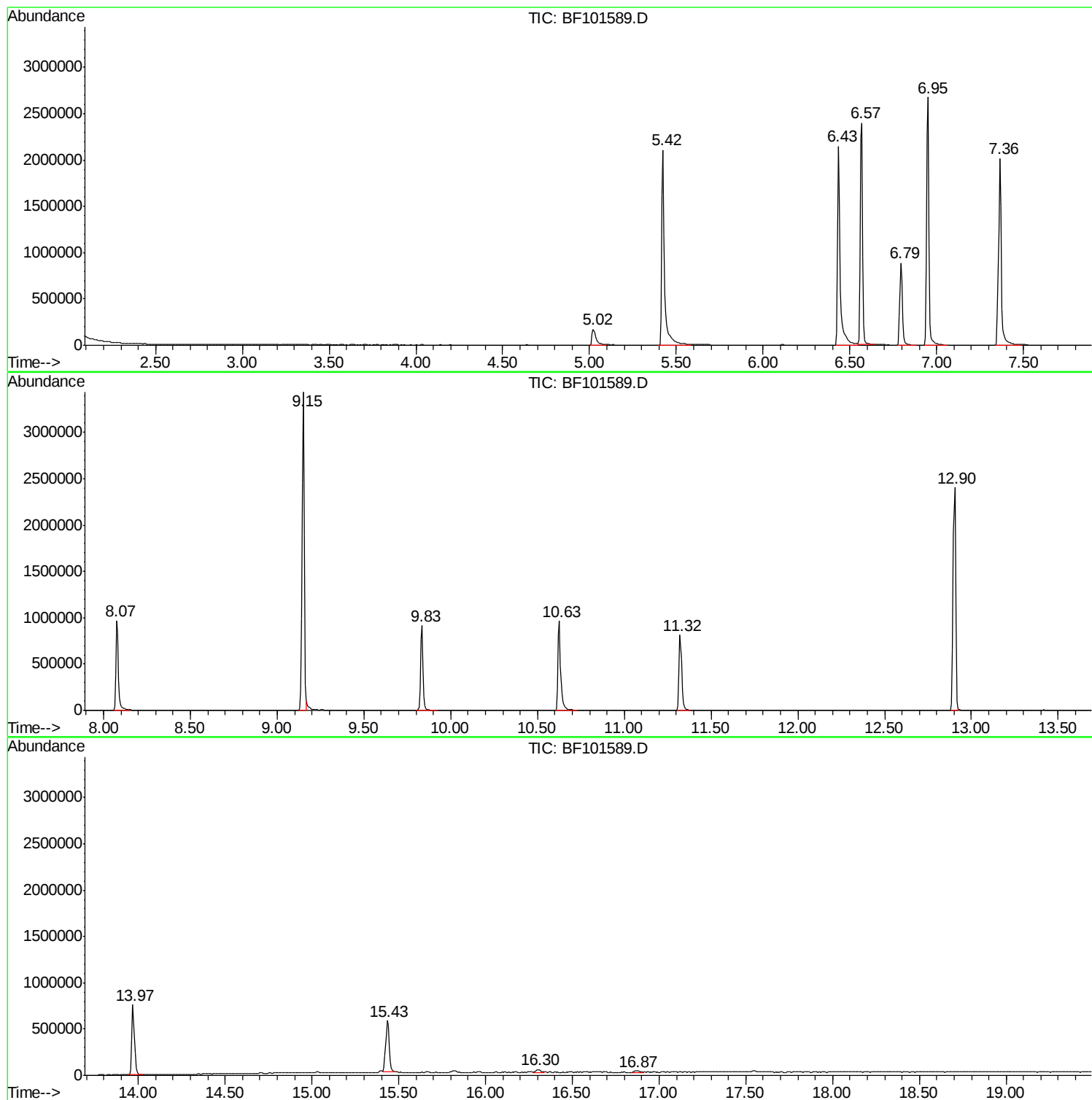
Sum of corrected areas: 24204154

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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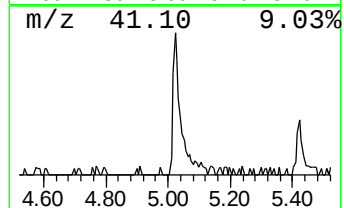
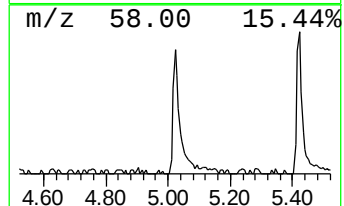
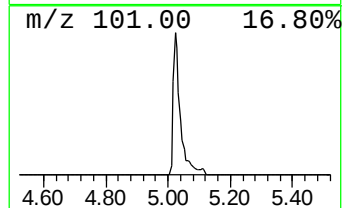
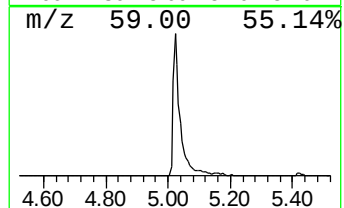
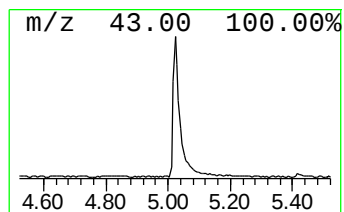
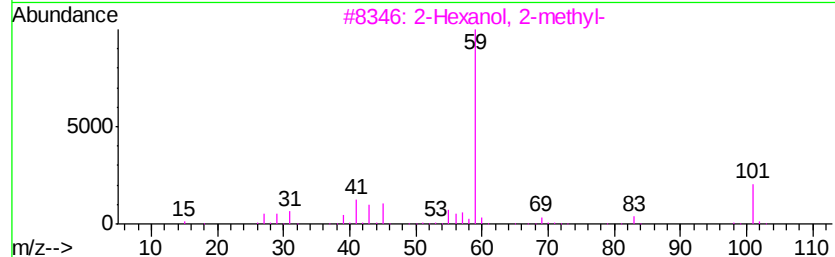
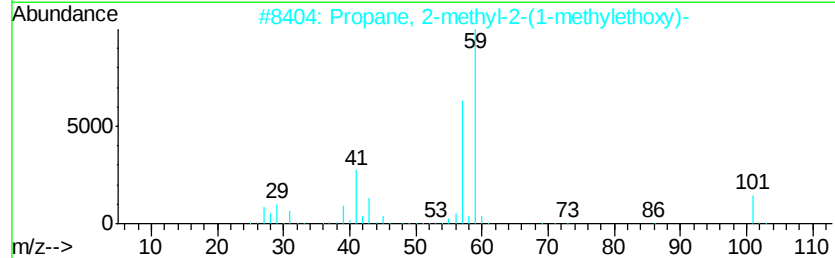
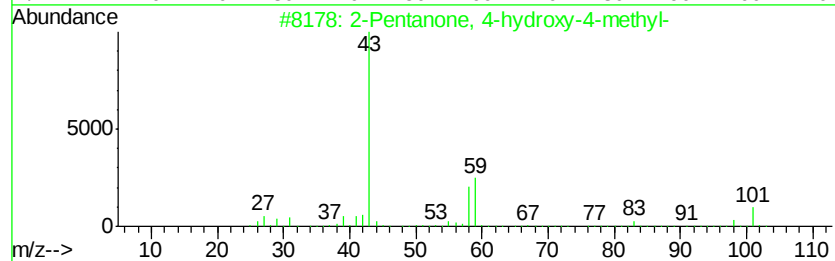
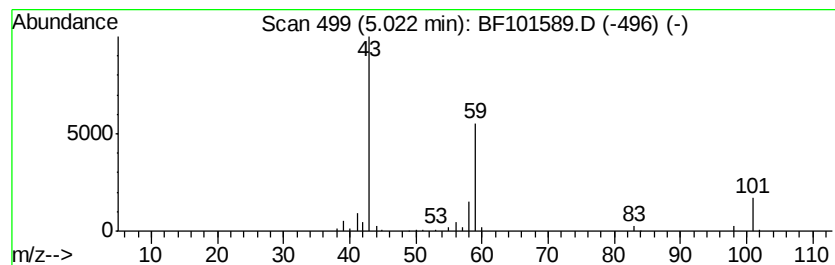
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	6.31 ng	263033	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	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O		017348-59-3	36
3	2-Hexanol, 2-methyl-	116	C7H16O		000625-23-0	33
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2		001116-98-9	25
5	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	10



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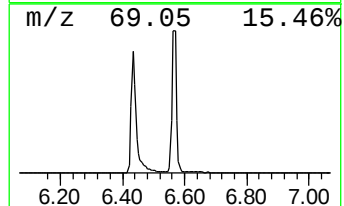
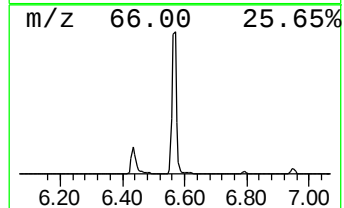
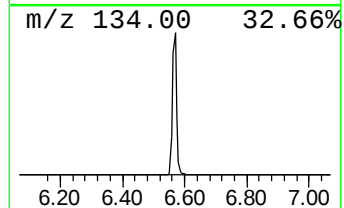
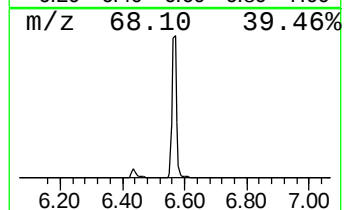
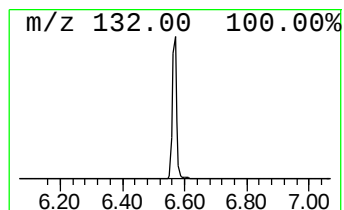
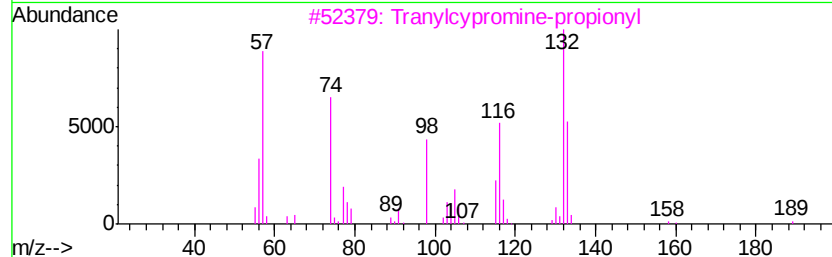
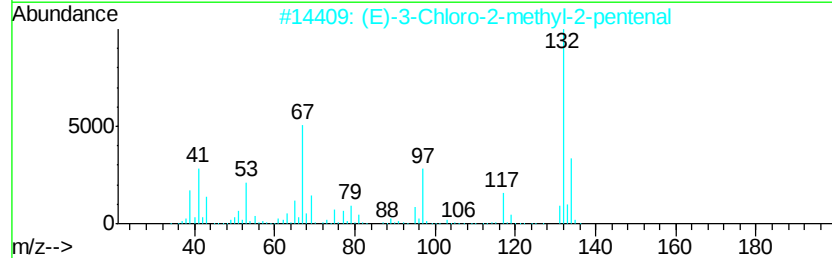
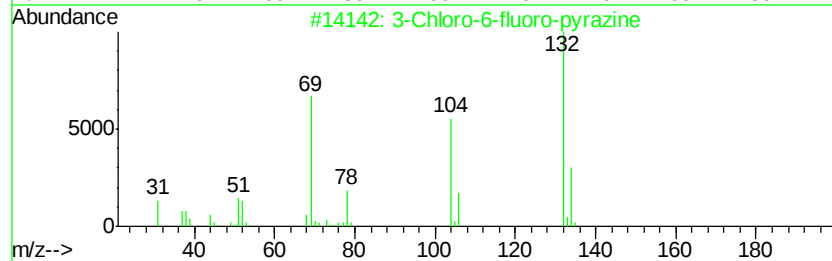
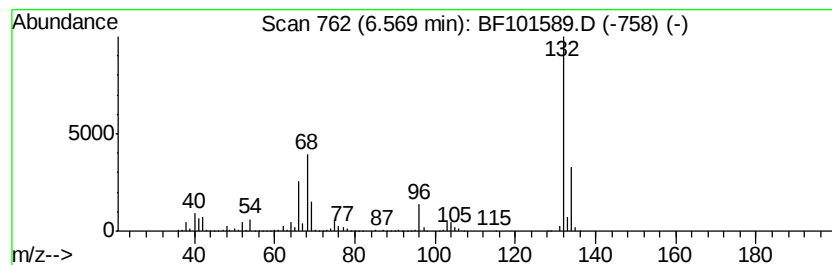
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Peak Number 2 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	57.45 ng	2394840	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
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...	5.02	6.3	ng	263033	1	6.79	833660	20.0
unknown6.57	6.57	57.5	ng	2394840	1	6.79	833660	20.0