

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102618\
 Data File : BG037594.D
 Acq On : 26 Oct 2018 23:43
 Operator : JU/SJ
 Sample : PB114204BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114204BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.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	5.180	318	322	331	rVB	38653	60065	1.66%	0.331%
2	5.644	393	401	413	rBV	837528	1361498	37.52%	7.499%
3	7.248	664	674	682	rBV	868283	1510447	41.62%	8.319%
4	7.630	729	739	752	rBV	797279	1419232	39.11%	7.817%
5	8.100	811	819	829	rVB	166830	299099	8.24%	1.647%
6	8.423	866	874	885	rBV	575370	1047265	28.86%	5.768%
7	9.275	1011	1019	1035	rBV	482108	911353	25.11%	5.020%
8	10.920	1292	1299	1308	rVB	242659	448395	12.36%	2.470%
9	13.359	1706	1714	1726	rBV	1396882	2275538	62.70%	12.533%
10	14.733	1941	1948	1955	rBV2	423246	696015	19.18%	3.834%
11	16.220	2194	2201	2227	rBV2	1060770	1711959	47.17%	9.429%
12	17.483	2408	2416	2424	rBV2	598123	933499	25.72%	5.142%
13	20.080	2852	2858	2870	rBV	2707248	3629008	100.00%	19.988%
14	21.766	3138	3145	3159	rBV2	548962	991615	27.32%	5.462%
15	25.098	3701	3712	3726	rBV3	247912	860827	23.72%	4.741%

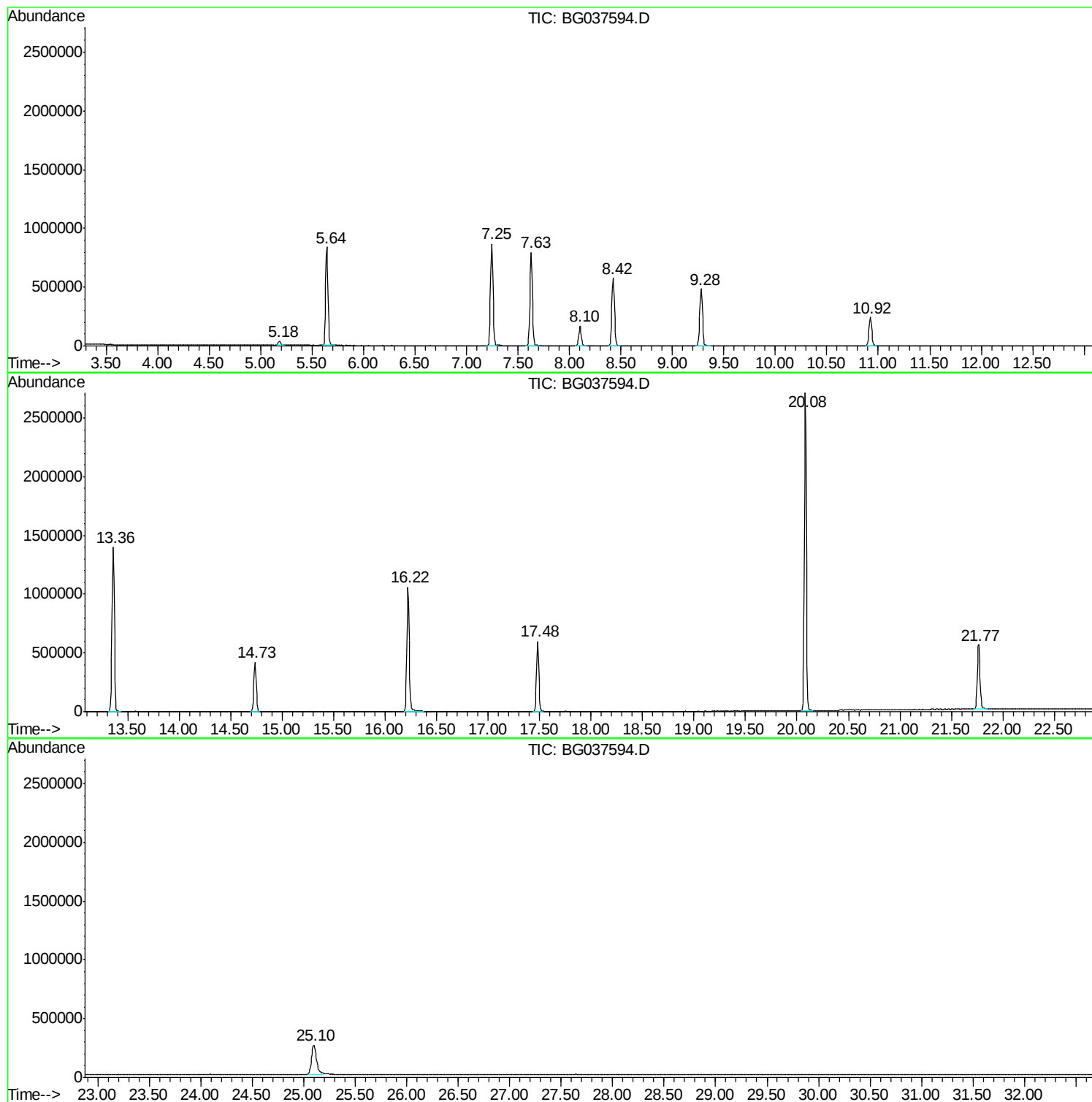
Sum of corrected areas: 18155815

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



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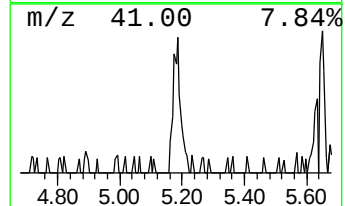
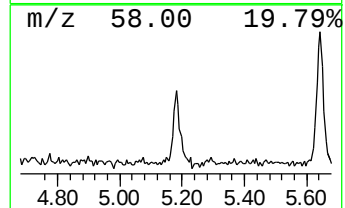
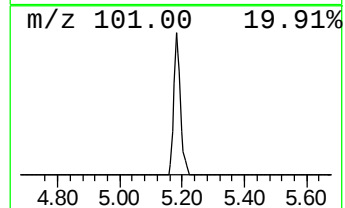
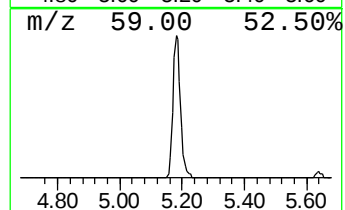
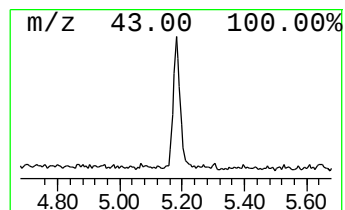
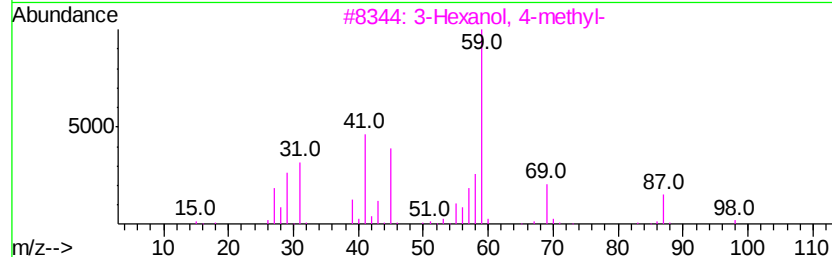
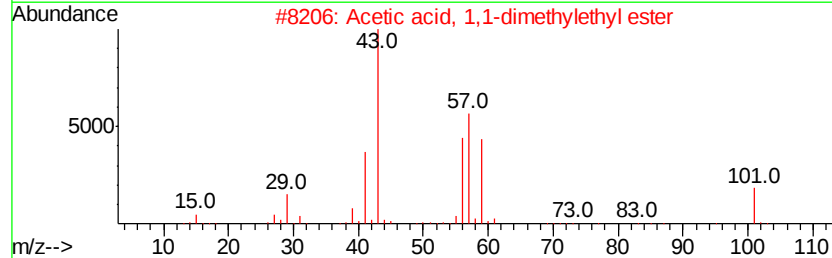
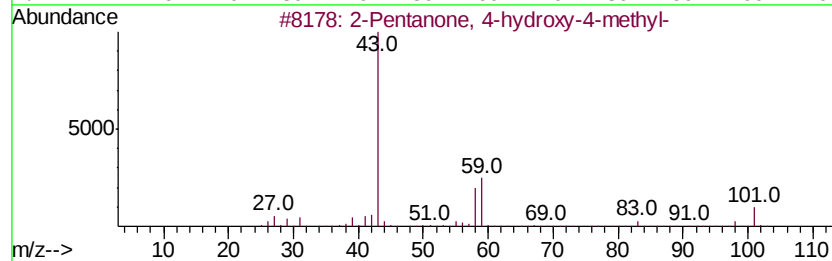
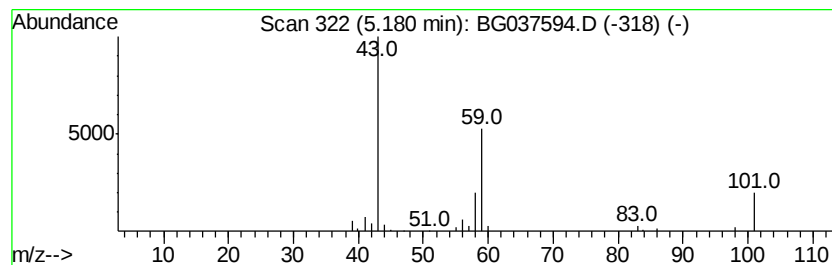
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.02 ng	60065	1,4-Dichlorobenzene-d4	8.10

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	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			3-Hexanol	102	C6H14O	000623-37-0	17
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12



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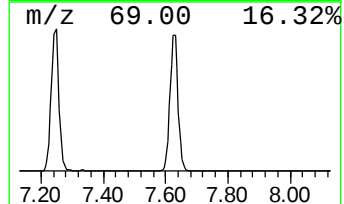
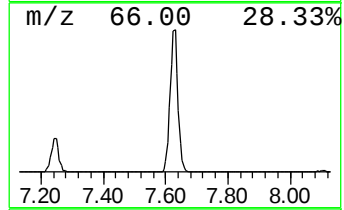
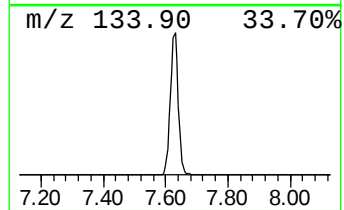
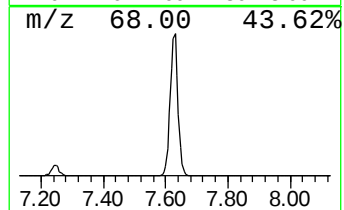
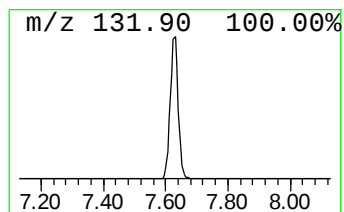
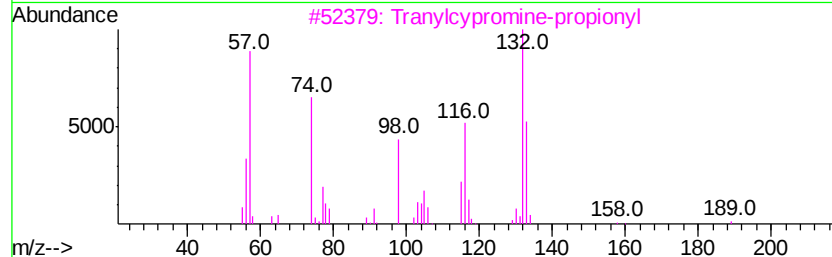
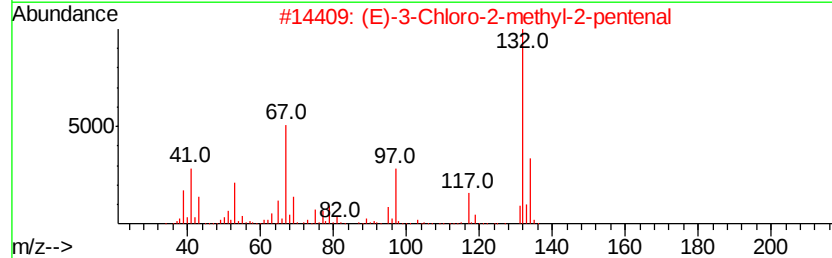
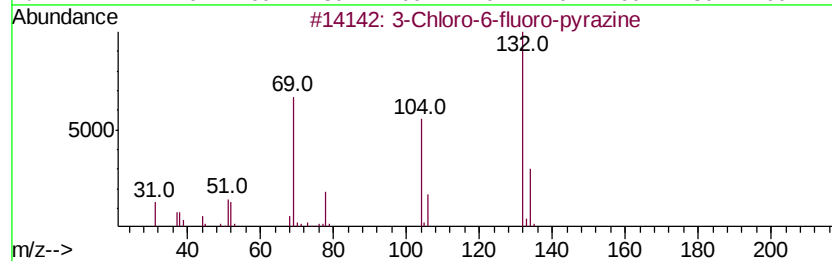
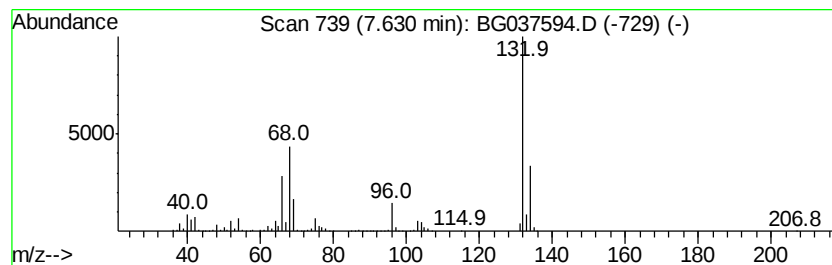
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	94.90 ng	1419230	1,4-Dichlorobenzene-d4	8.10

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	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	4.0	ng	60065	1	8.10	299099	20.0
unknown7.63	7.63	94.9	ng	1419230	1	8.10	299099	20.0