

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110218\
 Data File : BG037764.D
 Acq On : 2 Nov 2018 21:17
 Operator : JU/SJ
 Sample : PB114420BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114420BL

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	316	322	337	rVB	44936	79106	1.88%	0.360%
2	5.644	393	401	412	rBV	1022049	1727378	41.01%	7.854%
3	7.242	665	673	684	rBV	1090998	1956435	46.45%	8.896%
4	7.624	729	738	749	rBV	1034120	1832366	43.51%	8.332%
5	8.100	811	819	827	rBV	201764	351534	8.35%	1.598%
6	8.423	866	874	885	rBV	763219	1385858	32.90%	6.302%
7	9.275	1011	1019	1030	rBV	656609	1203093	28.56%	5.470%
8	10.920	1291	1299	1310	rBV	282689	545089	12.94%	2.479%
9	13.353	1706	1713	1726	rBV	1907573	3127927	74.27%	14.223%
10	14.734	1941	1948	1960	rBV2	502737	831019	19.73%	3.779%
11	16.220	2194	2201	2223	rBV	1274538	2052527	48.73%	9.333%
12	17.477	2406	2415	2427	rBV2	606017	974998	23.15%	4.433%
13	20.080	2851	2858	2873	rBV	3048827	4211776	100.00%	19.151%
14	21.767	3137	3145	3165	rBV	513058	1015358	24.11%	4.617%
15	25.104	3702	3713	3732	rBV3	186357	698038	16.57%	3.174%

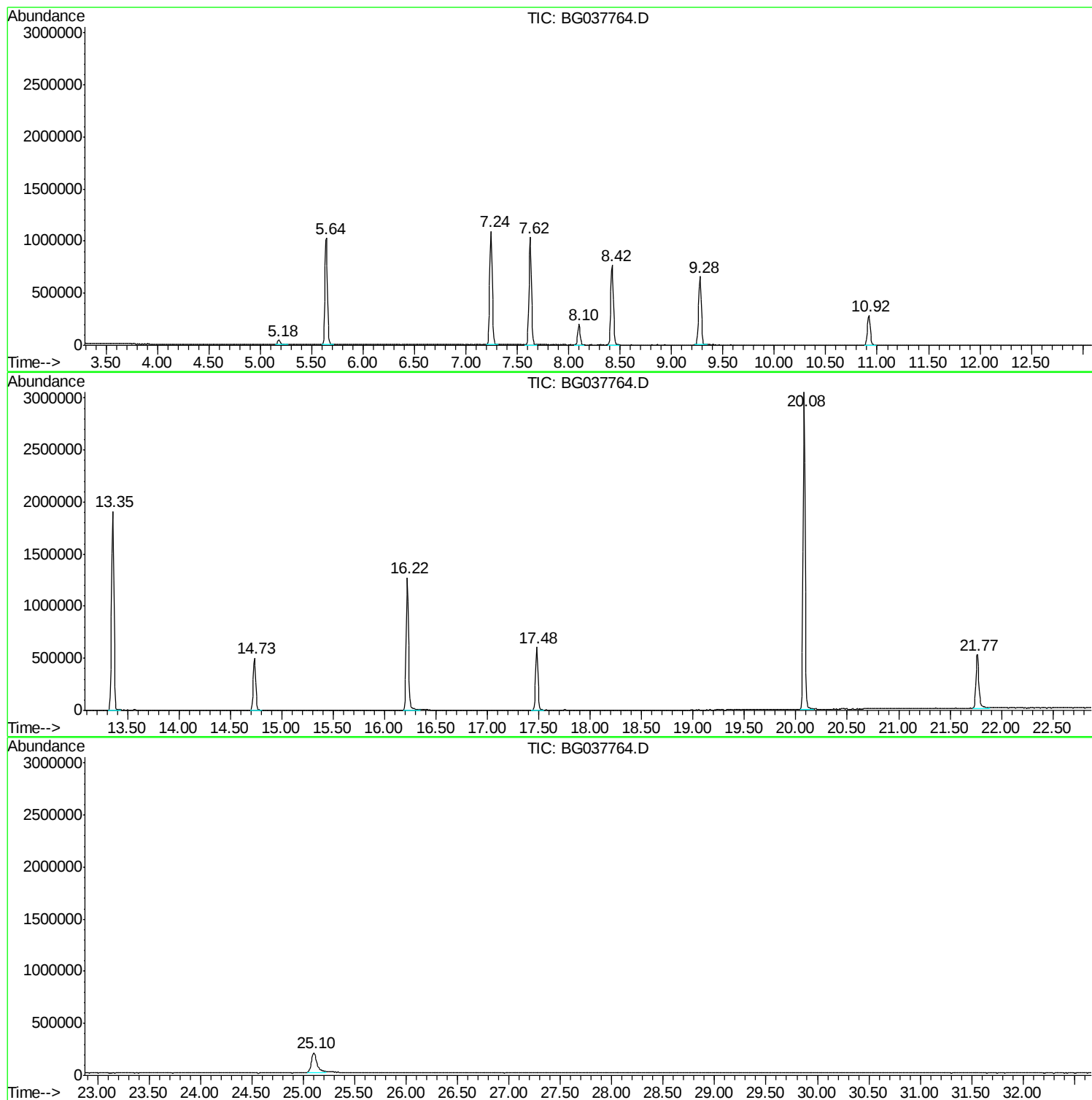
Sum of corrected areas: 21992502

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



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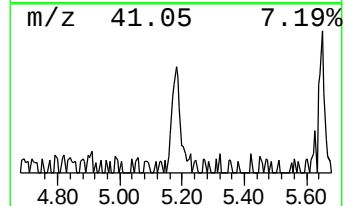
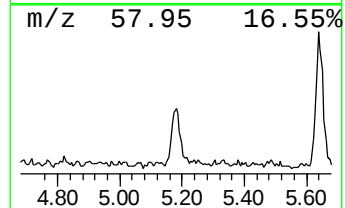
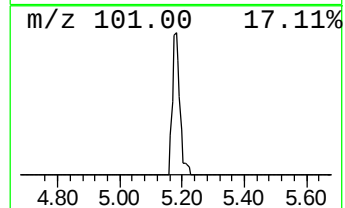
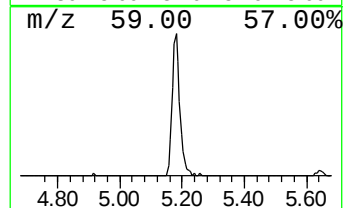
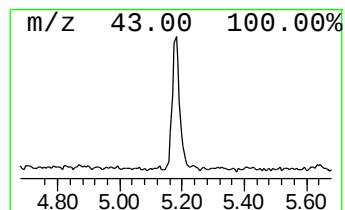
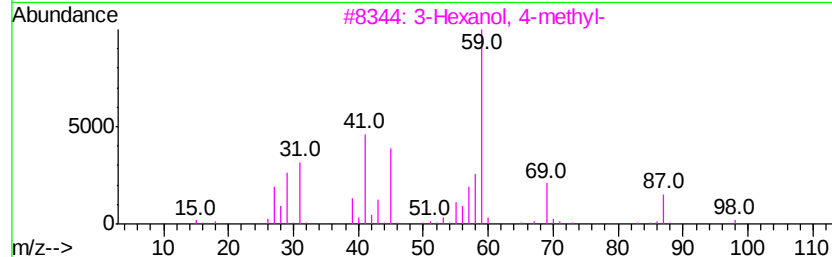
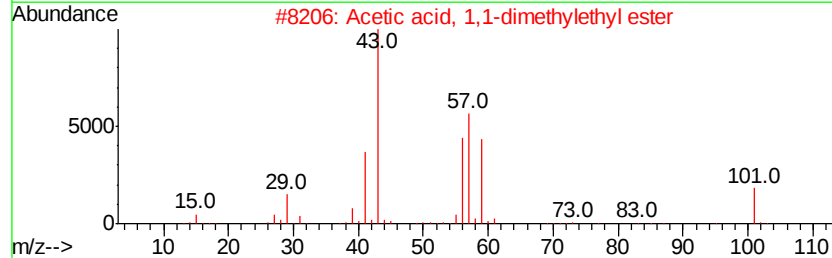
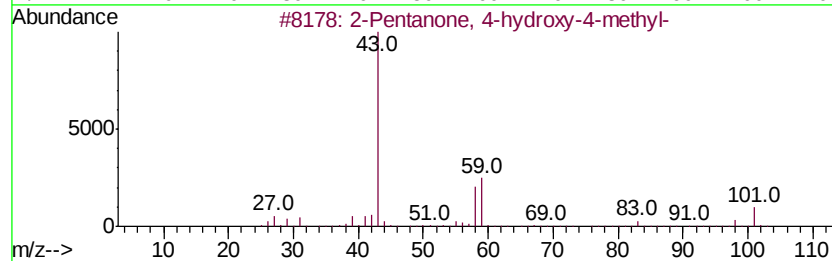
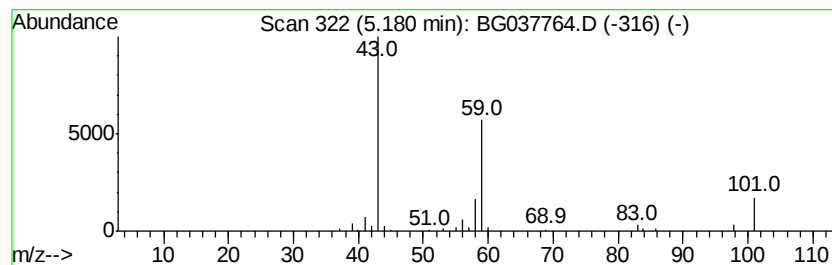
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.50 ng	79106	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	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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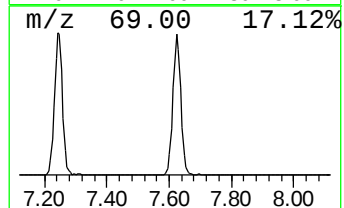
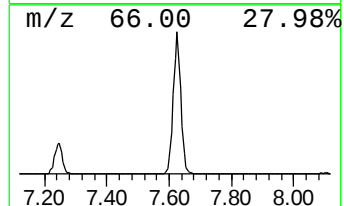
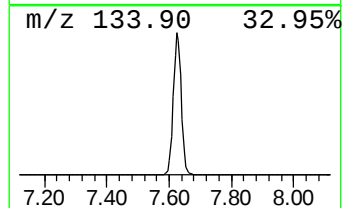
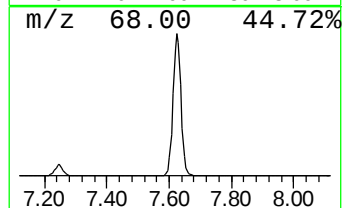
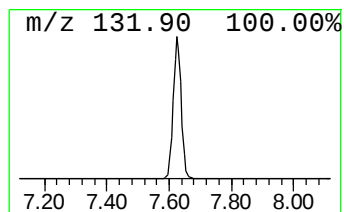
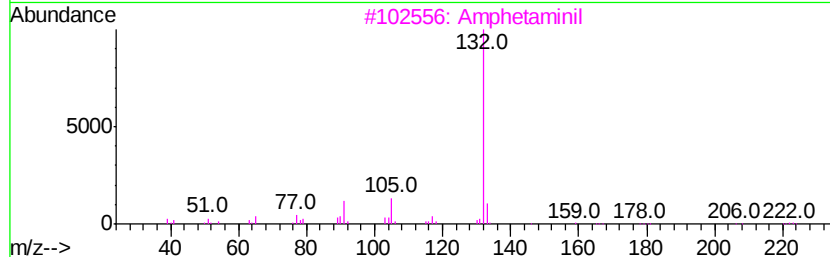
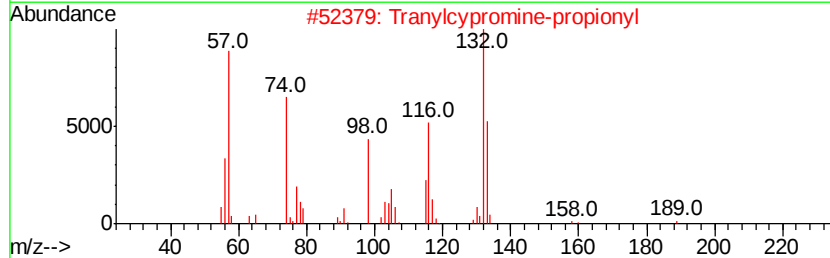
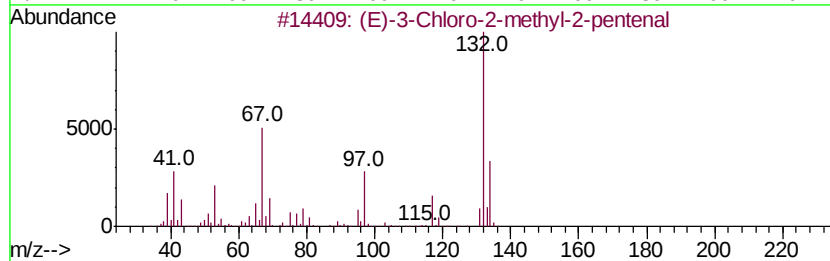
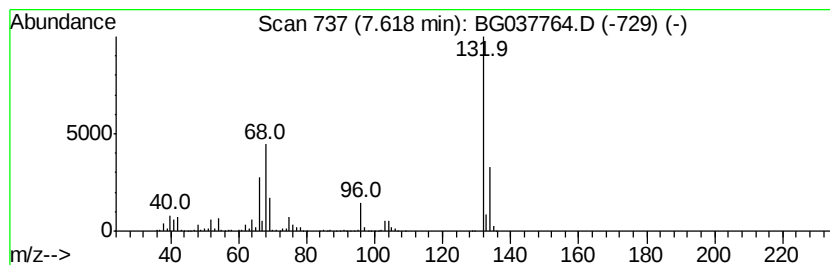
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	104.25 ng	1832370	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Amphetaminil	250	C17H18N2	017590-01-1	10
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	4.5	ng	79106	1	8.10	351534	20.0
unknown7.62	7.62	104.3	ng	1832370	1	8.10	351534	20.0