

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG110618\
 Data File : BG037841.D
 Acq On : 6 Nov 2018 16:47
 Operator : JU/SJ
 Sample : PB114487BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114487BL

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	5.176	316	321	330	rVB	36299	58897	1.71%	0.328%
2	5.634	392	399	413	rBV	822114	1392028	40.32%	7.755%
3	7.238	663	672	683	rBV	901835	1585229	45.91%	8.831%
4	7.620	729	737	748	rBV	842535	1502977	43.53%	8.373%
5	8.096	810	818	825	rBV	173584	302433	8.76%	1.685%
6	8.419	865	873	881	rBV	614168	1127258	32.65%	6.280%
7	9.271	1009	1018	1033	rBV	519017	993930	28.79%	5.537%
8	10.916	1290	1298	1305	rBV	253472	480818	13.93%	2.678%
9	13.349	1705	1712	1727	rBV	1575306	2561328	74.18%	14.268%
10	14.729	1939	1947	1962	rBV2	436420	727019	21.06%	4.050%
11	16.216	2193	2200	2215	rBV2	895127	1496406	43.34%	8.336%
12	17.473	2408	2414	2428	rBV2	543828	848534	24.58%	4.727%
13	20.076	2851	2857	2875	rBV	2488074	3452765	100.00%	19.234%
14	21.762	3137	3144	3157	rBV	422788	846347	24.51%	4.715%
15	25.094	3700	3711	3728	rBV2	143698	575224	16.66%	3.204%

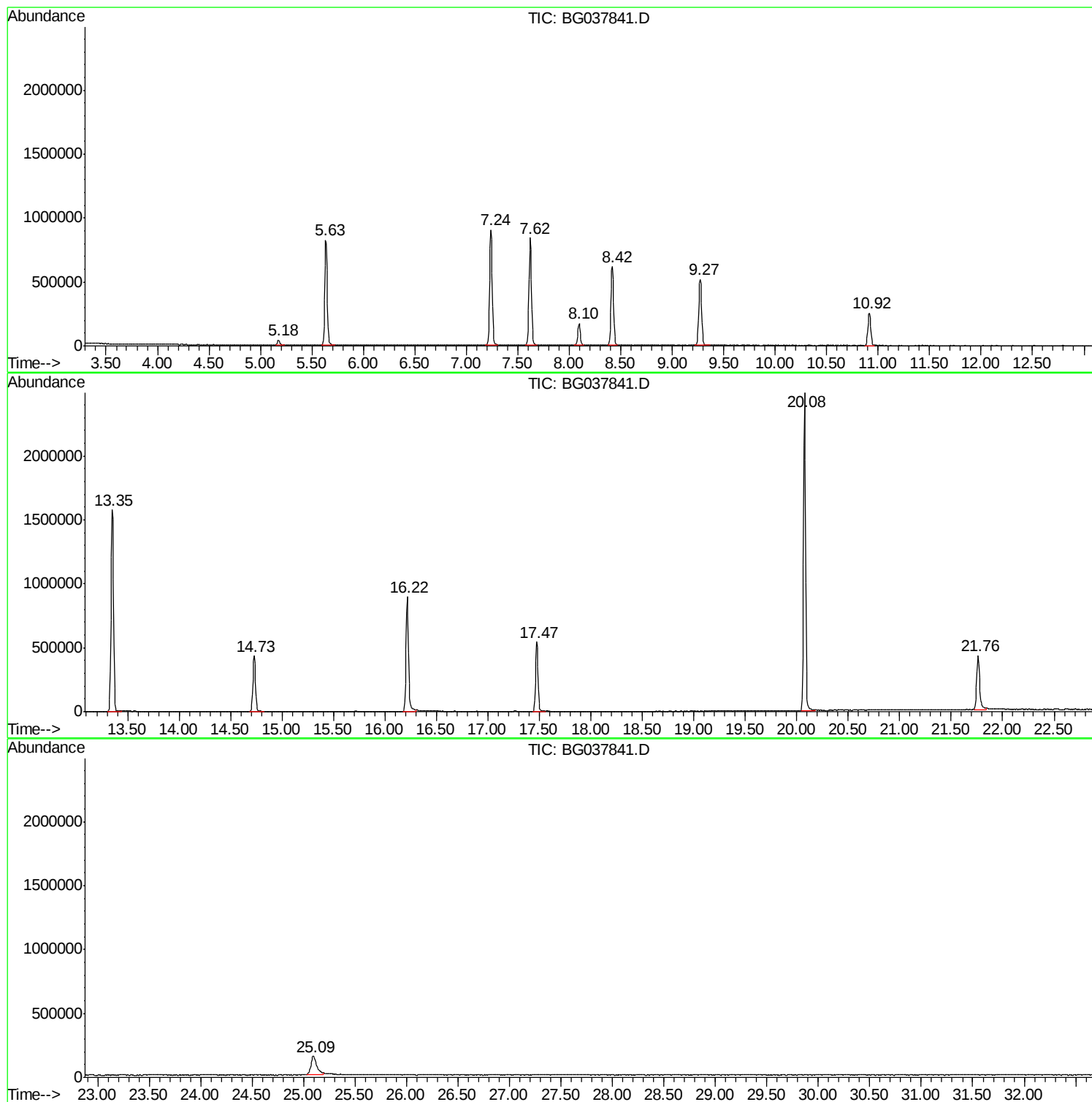
Sum of corrected areas: 17951193

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



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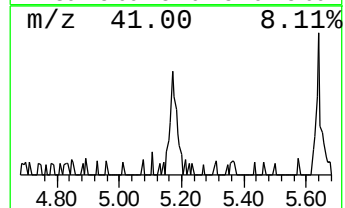
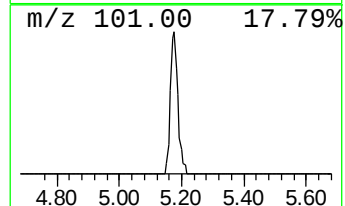
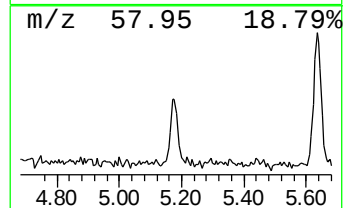
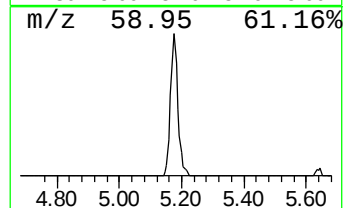
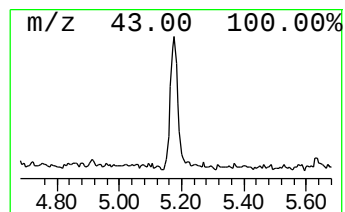
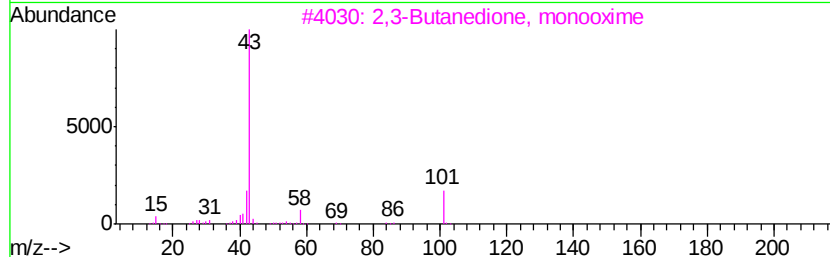
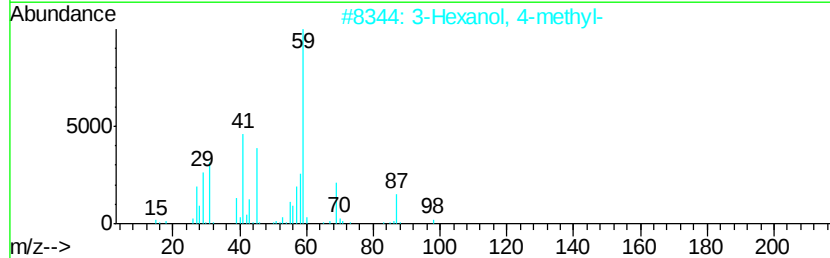
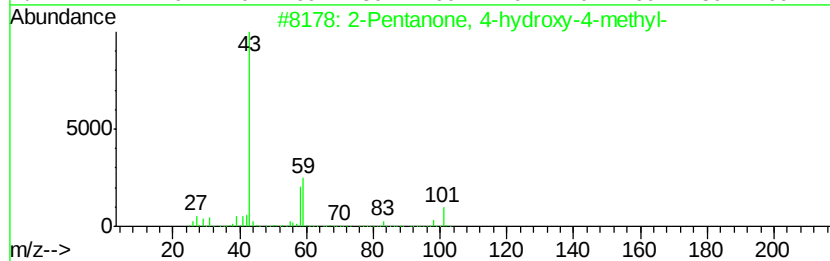
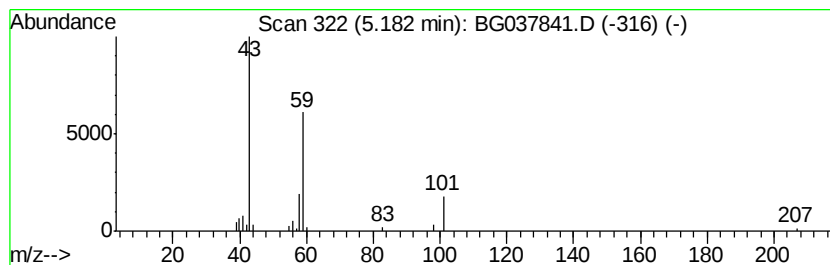
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.89 ng	58897	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	53
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Acetone	58	C3H6O	000067-64-1	9



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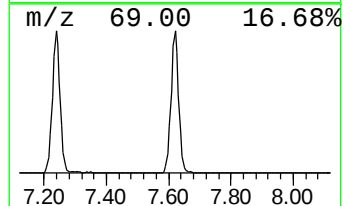
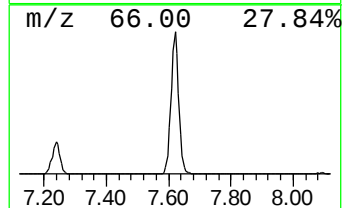
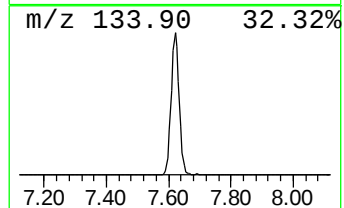
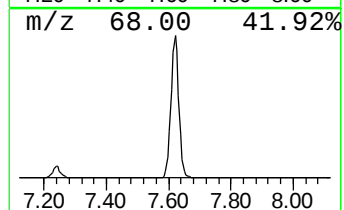
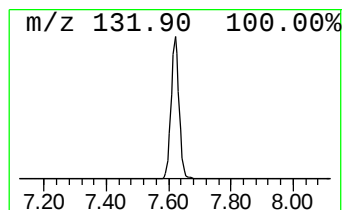
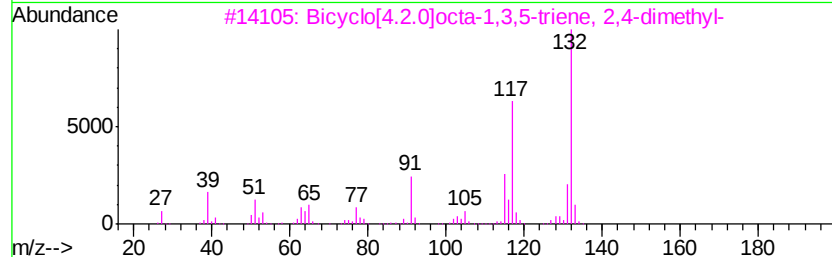
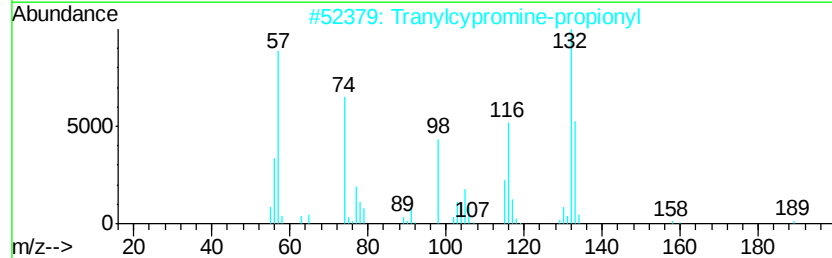
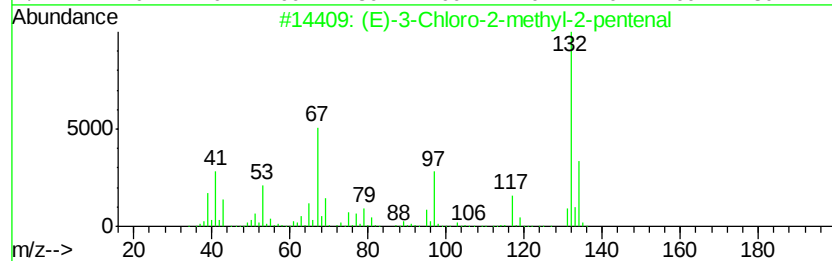
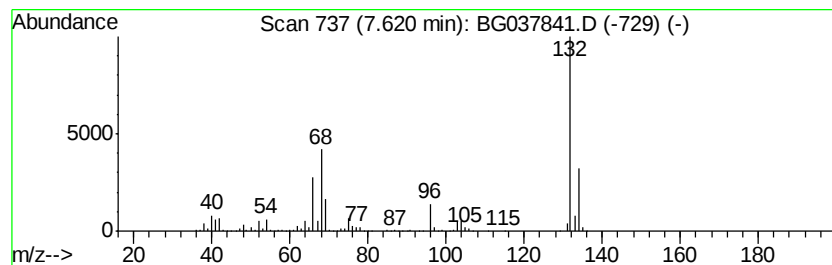
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	99.39 ng	1502980	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	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.9	ng	58897	1	8.10	302433	20.0
unknown7.62	7.62	99.4	ng	1502980	1	8.10	302433	20.0