

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110518\
 Data File : BG037823.D
 Acq On : 6 Nov 2018 4:11
 Operator : JU/SJ
 Sample : PB114495BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114495BL

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	314	321	329	rBV	51976	87702	1.85%	0.355%
2	5.640	392	400	414	rBV	1194026	1984580	41.77%	8.024%
3	7.244	664	673	685	rBV	1251788	2240470	47.16%	9.058%
4	7.620	729	737	753	rBV	1172048	2117941	44.58%	8.563%
5	8.096	810	818	825	rBV	211067	378299	7.96%	1.529%
6	8.419	865	873	881	rBV	893614	1596311	33.60%	6.454%
7	9.271	1010	1018	1030	rBV	749697	1410077	29.68%	5.701%
8	10.916	1290	1298	1307	rBV	319902	601965	12.67%	2.434%
9	13.349	1705	1712	1722	rBV	2174853	3556652	74.86%	14.380%
10	14.730	1939	1947	1959	rBV2	533426	898173	18.90%	3.631%
11	16.216	2193	2200	2219	rBV	1378603	2214727	46.61%	8.954%
12	17.473	2407	2414	2425	rBV2	658076	1047164	22.04%	4.234%
13	20.076	2851	2857	2872	rBV	3347132	4751257	100.00%	19.209%
14	21.757	3136	3143	3160	rBV2	535276	1055303	22.21%	4.267%
15	25.088	3699	3710	3725	rBV2	209797	793524	16.70%	3.208%

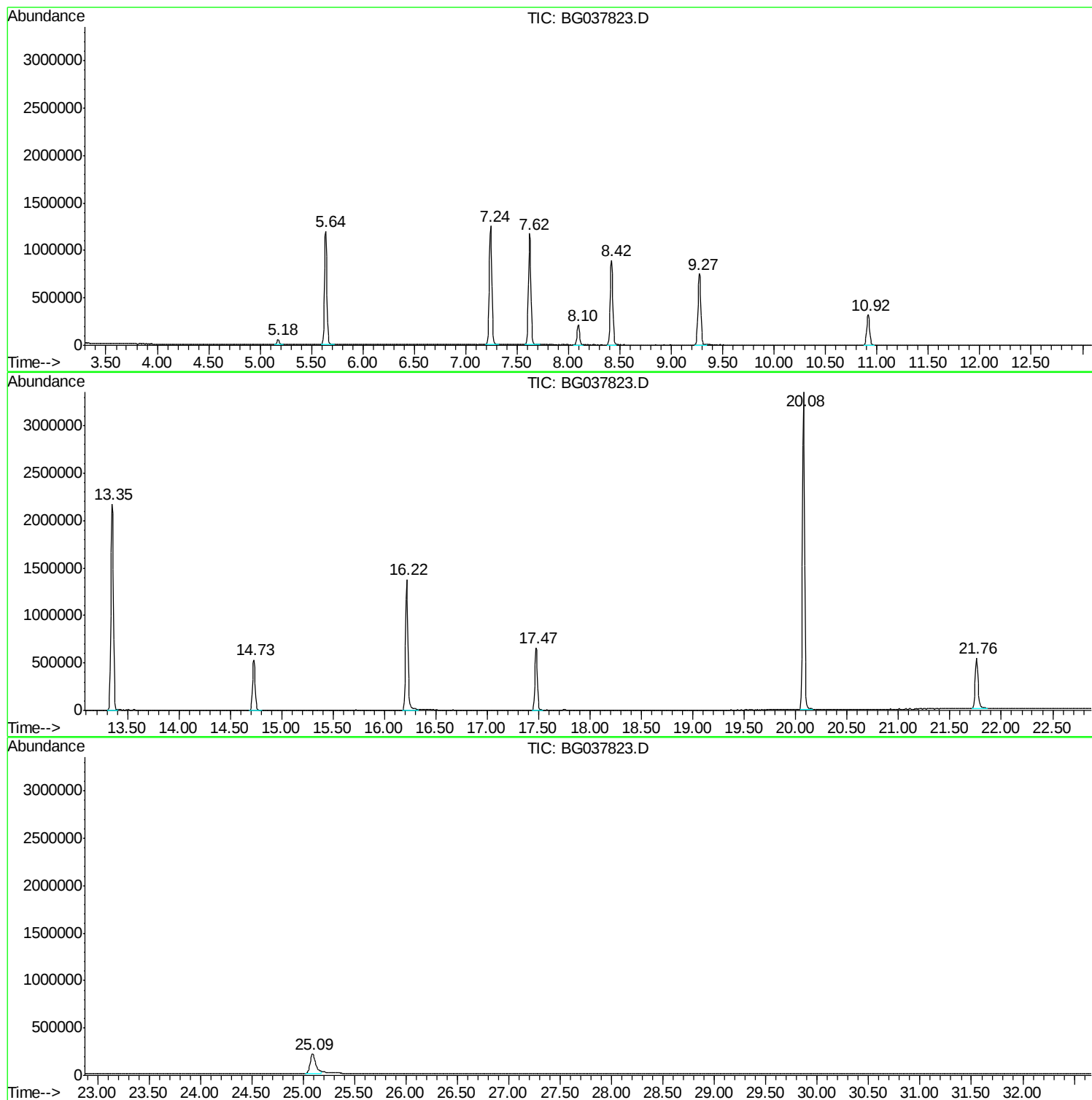
Sum of corrected areas: 24734145

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



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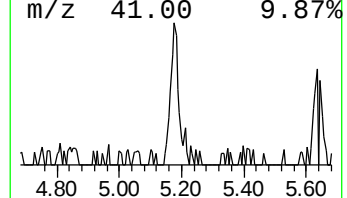
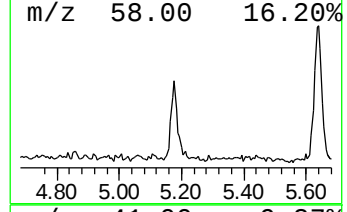
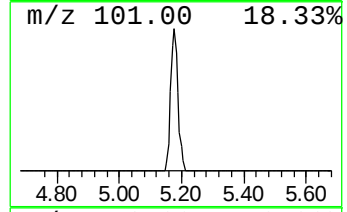
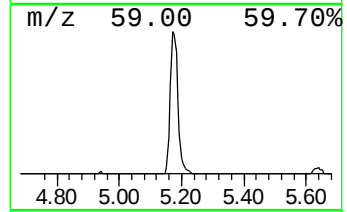
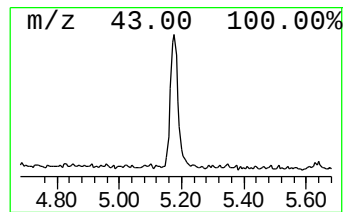
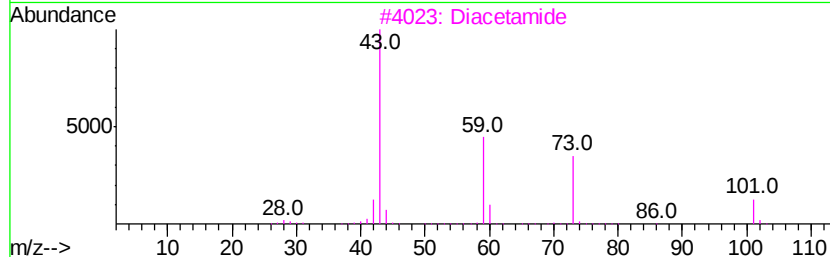
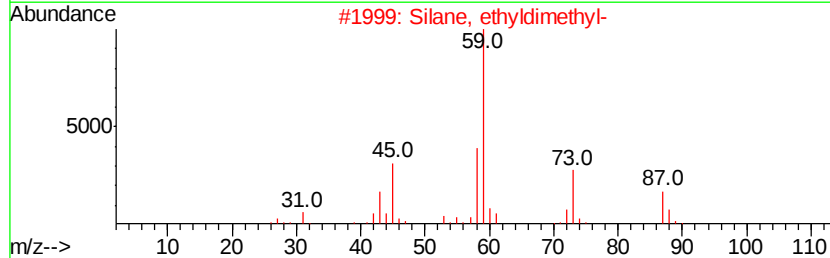
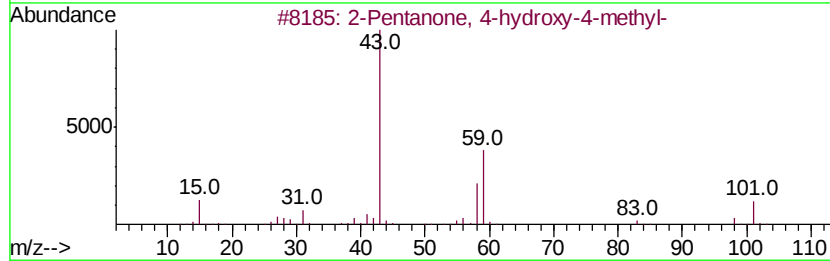
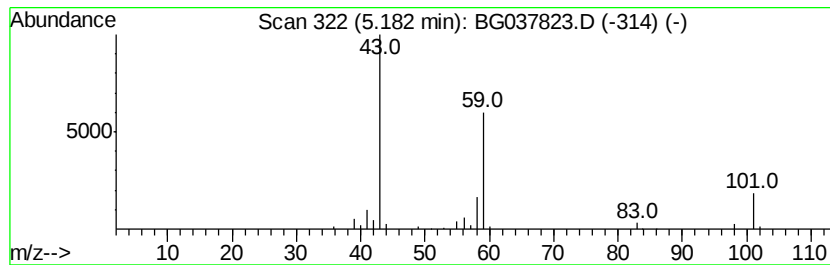
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.64 ng	87702	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	Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	28
3	Diacetamide	101	C4H7NO2	000625-77-4	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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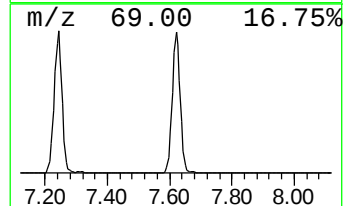
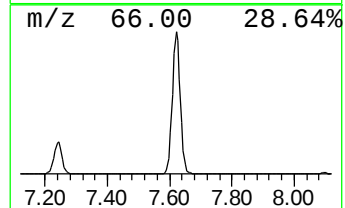
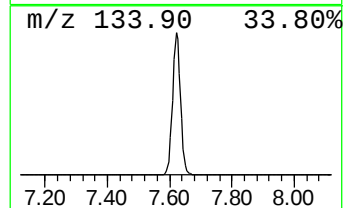
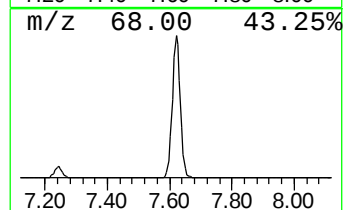
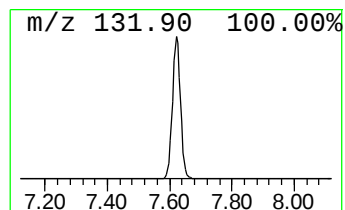
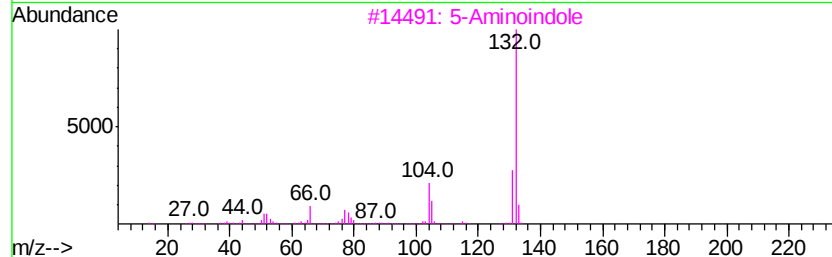
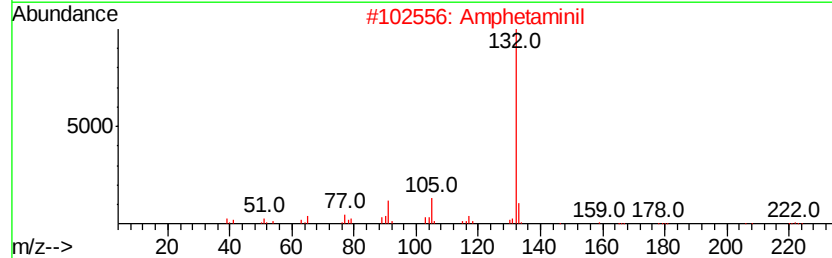
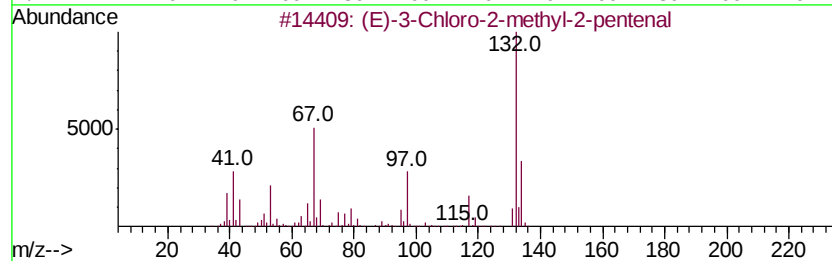
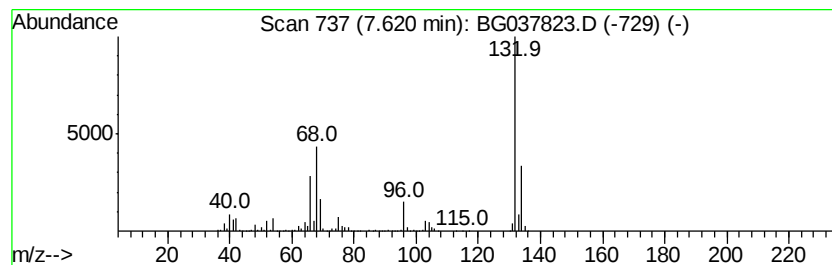
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	111.97 ng	2117940	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2			Amphetaminil	250	C17H18N2	017590-01-1	12
3			5-Aminoindole	132	C8H8N2	005192-03-0	12
4			Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	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...	5.18	4.6	ng	87702	1	8.10	378299	20.0
unknown7.62	7.62	112.0	ng	2117940	1	8.10	378299	20.0