

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030648.D
 Acq On : 1 Nov 2017 23:54
 Operator : SJ/JU
 Sample : I6204-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102617-TIDO-E-D-S

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_G\METHODS\8270-BG101617.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	5.242	325	332	341	rBV	102562	161088	2.57%	0.655%
2	5.718	405	413	423	rBV	457592	756985	12.08%	3.079%
3	7.316	677	685	696	rBV	307525	588740	9.39%	2.395%
4	7.692	739	749	758	rBV	772480	1358007	21.67%	5.523%
5	8.156	822	828	836	rVB	215047	370819	5.92%	1.508%
6	8.485	876	884	892	rBV	1016213	1796217	28.66%	7.306%
7	9.325	1017	1027	1037	rBV	811372	1496716	23.88%	6.088%
8	10.971	1299	1307	1317	rBV	324051	593987	9.48%	2.416%
9	13.397	1713	1720	1729	rBV	2619881	4140631	66.06%	16.841%
10	14.214	1852	1859	1868	rBV	43970	70162	1.12%	0.285%
11	14.772	1947	1954	1965	rBV2	595008	982124	15.67%	3.995%
12	16.253	2198	2206	2228	rBV	1315043	2092456	33.39%	8.511%
13	17.510	2414	2420	2431	rBV	732583	1148233	18.32%	4.670%
14	20.107	2855	2862	2875	rBV	4566504	6267577	100.00%	25.492%
15	21.781	3140	3147	3155	rBV2	821480	1380182	22.02%	5.614%
16	25.089	3699	3710	3727	rBV2	471237	1382685	22.06%	5.624%

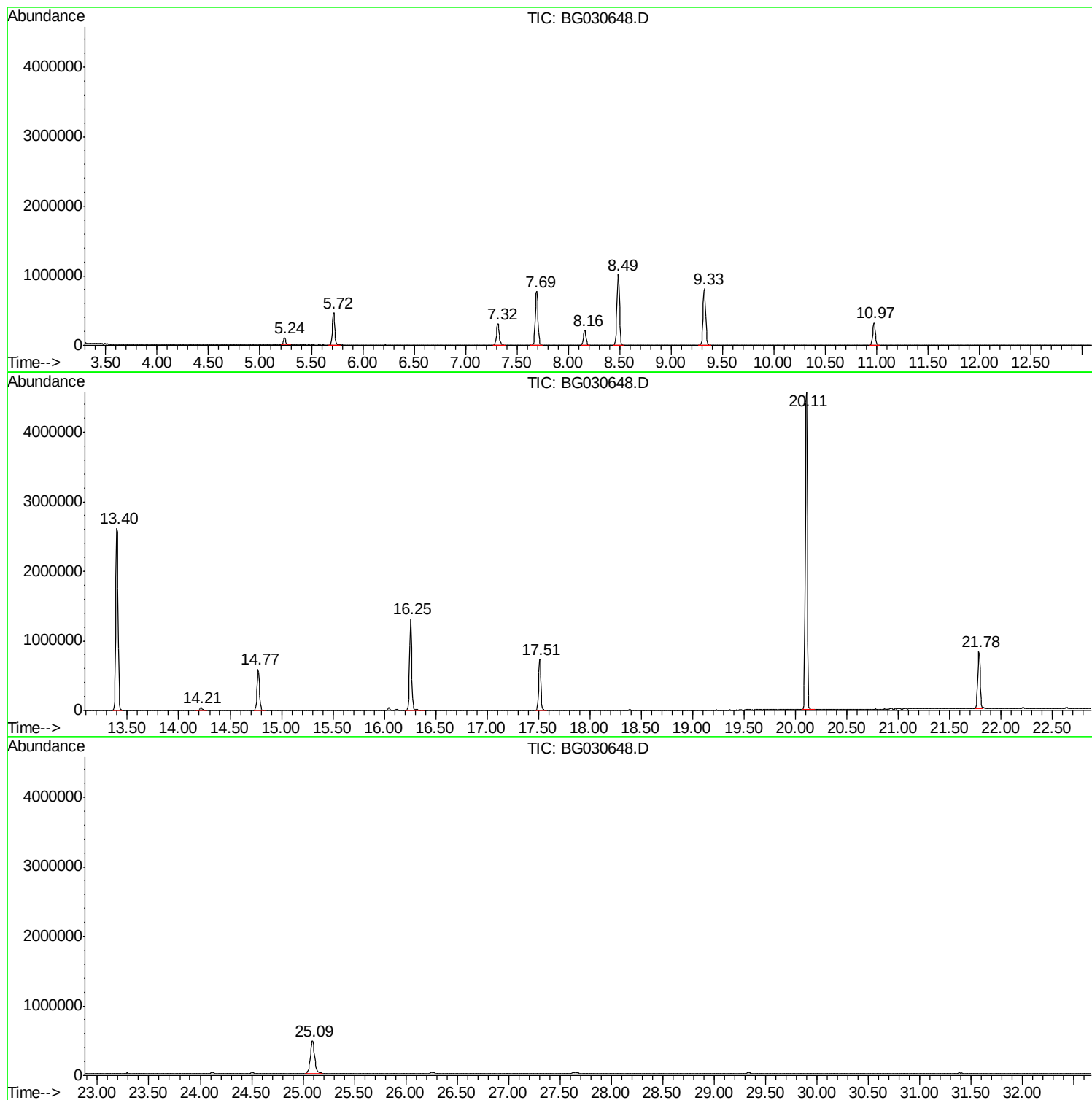
Sum of corrected areas: 24586609

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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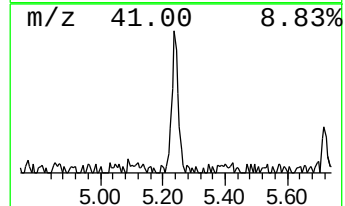
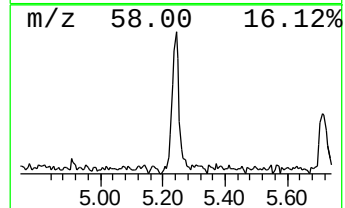
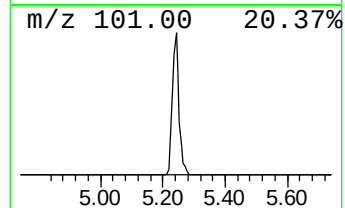
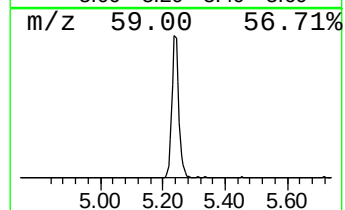
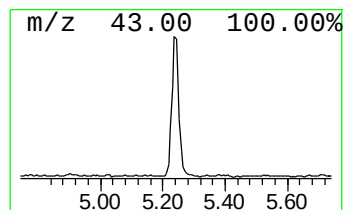
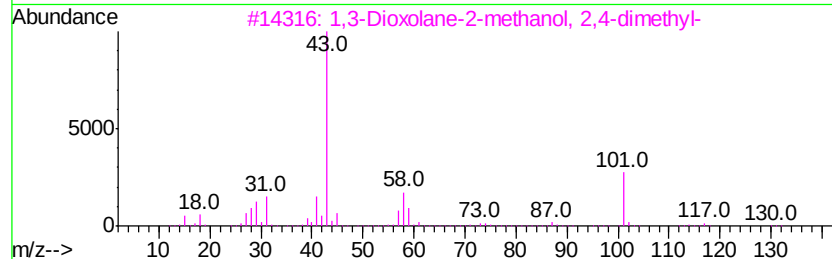
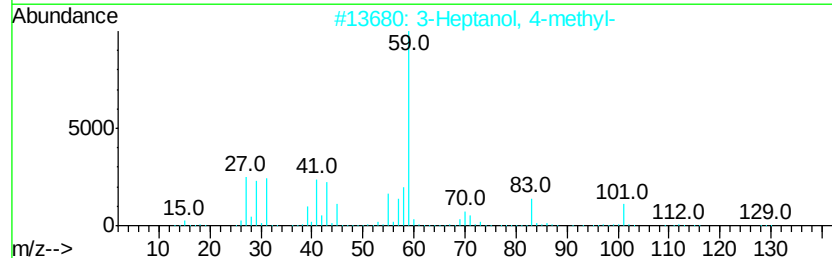
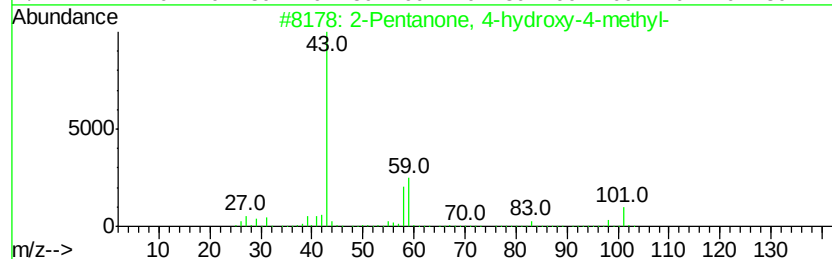
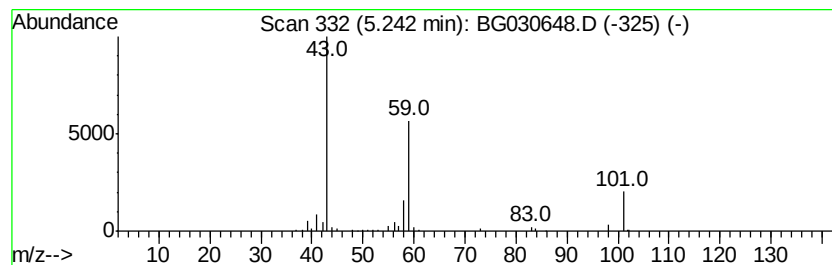
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	8.69 ng	161088	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	28
3			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	25
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	16
5			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	12



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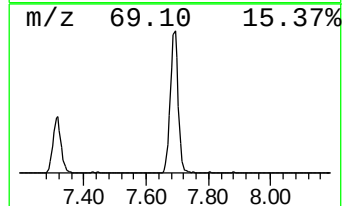
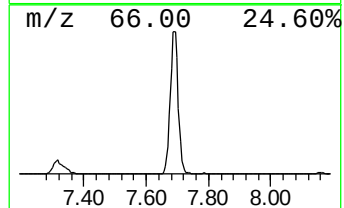
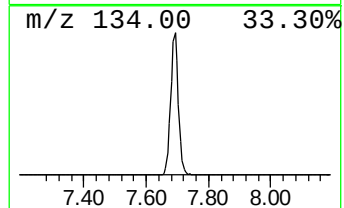
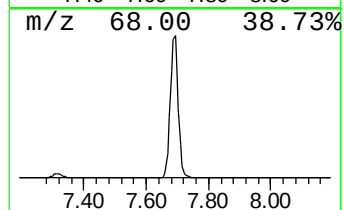
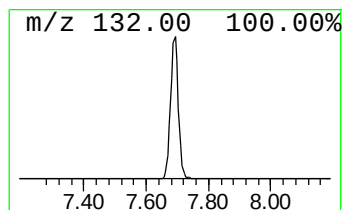
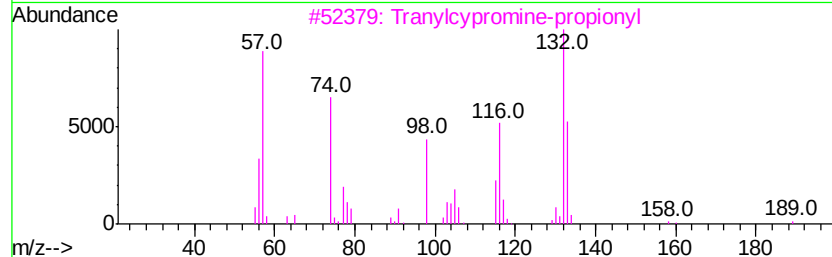
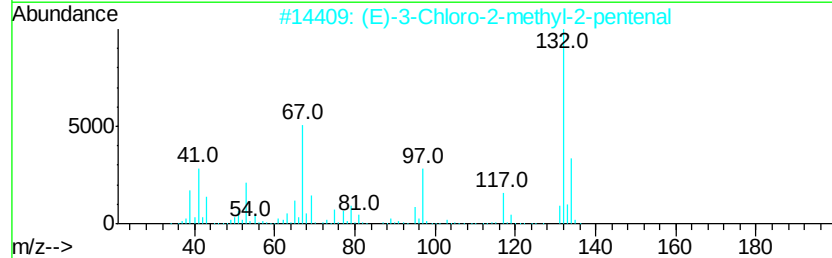
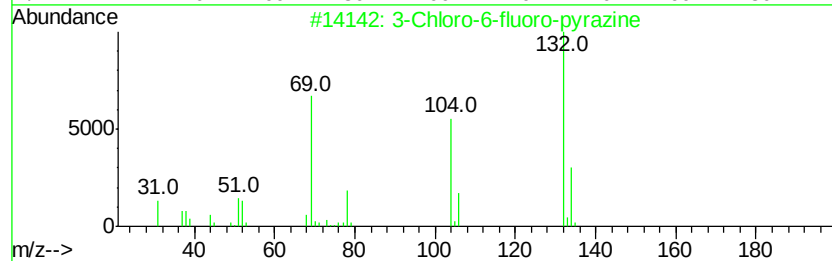
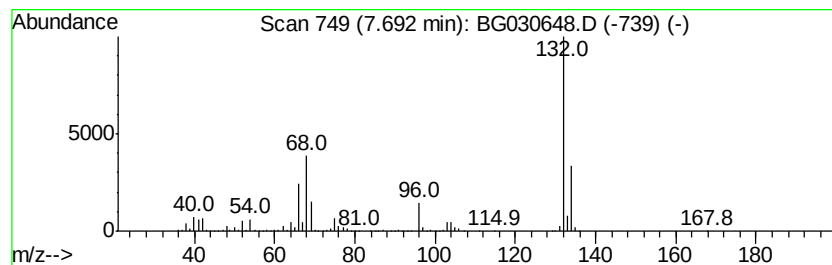
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Peak Number 2 unknown7.69 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	73.24 ng	1358010	1,4-Dichlorobenzene-d4	8.16

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	10
4		Amphetaminil	250	C17H18N2	017590-01-1	10
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	8.7	ng	161088	1	8.16	370819	20.0
unknown7.69	7.69	73.2	ng	1358010	1	8.16	370819	20.0