

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102817\
 Data File : BG030542.D
 Acq On : 29 Oct 2017 1:21
 Operator : SJ/JU
 Sample : PB103762BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103762BL

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	5.239	325	332	340	rBV	175386	266213	5.63%	1.185%
2	5.715	405	413	427	rBV	998000	1619117	34.27%	7.209%
3	7.319	678	686	697	rBV	981823	1702260	36.03%	7.579%
4	7.695	742	750	760	rBV	989512	1687980	35.73%	7.515%
5	8.165	823	830	837	rBV	191966	335442	7.10%	1.494%
6	8.488	878	885	899	rBV	738615	1307966	27.68%	5.824%
7	9.329	1020	1028	1037	rBV	574547	1007571	21.33%	4.486%
8	10.980	1301	1309	1320	rVB	268346	481586	10.19%	2.144%
9	13.406	1715	1722	1736	rBV	1739020	2711156	57.38%	12.071%
10	14.781	1949	1956	1968	rVB2	465514	738895	15.64%	3.290%
11	16.262	2201	2208	2231	rBV	1529306	2388005	50.54%	10.632%
12	17.519	2415	2422	2433	rBV2	685474	1030781	21.82%	4.589%
13	20.110	2857	2863	2873	rBV	3656905	4724561	100.00%	21.035%
14	21.790	3142	3149	3159	rBV	752147	1243795	26.33%	5.538%
15	25.098	3700	3712	3724	rBV2	415391	1214785	25.71%	5.409%

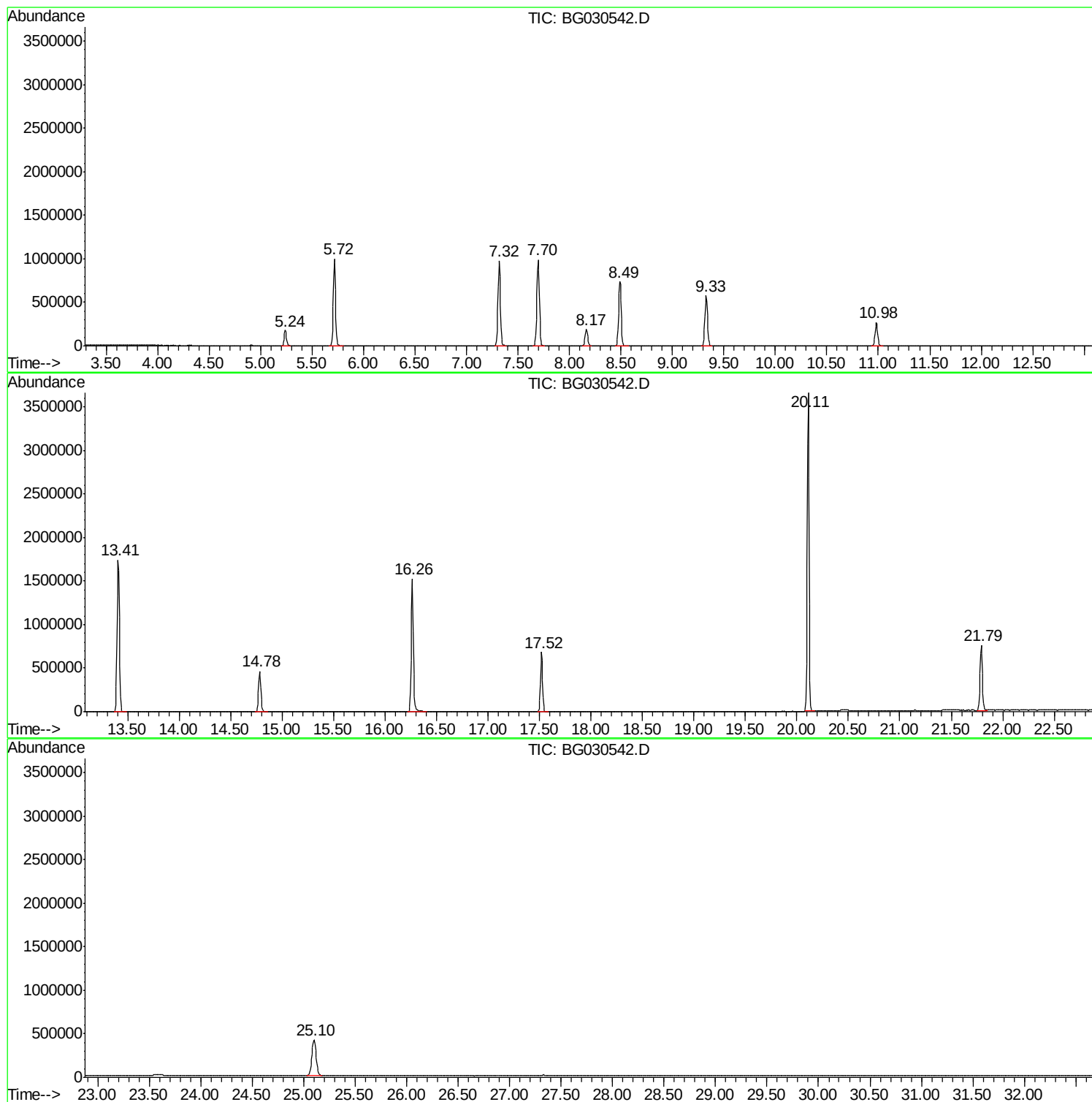
Sum of corrected areas: 22460113

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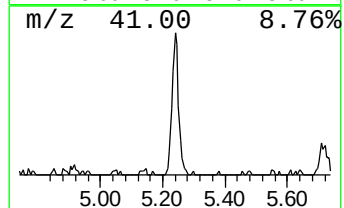
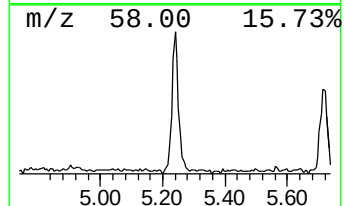
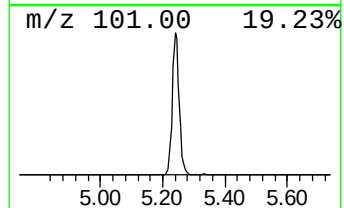
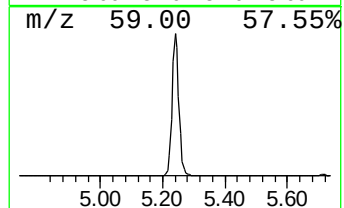
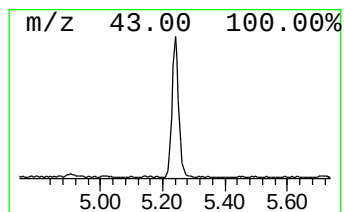
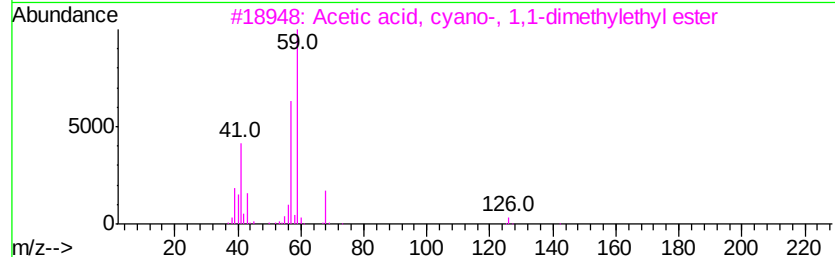
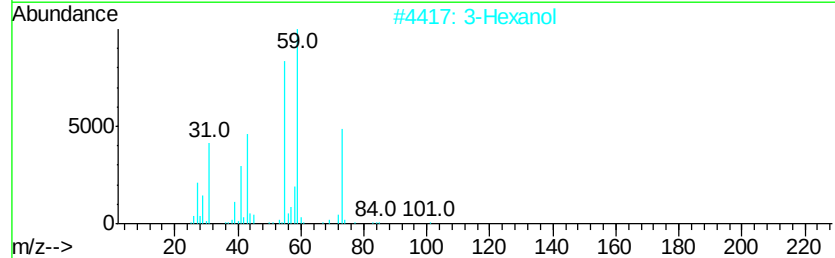
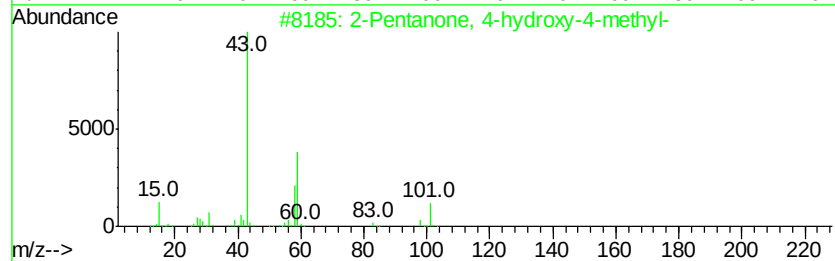
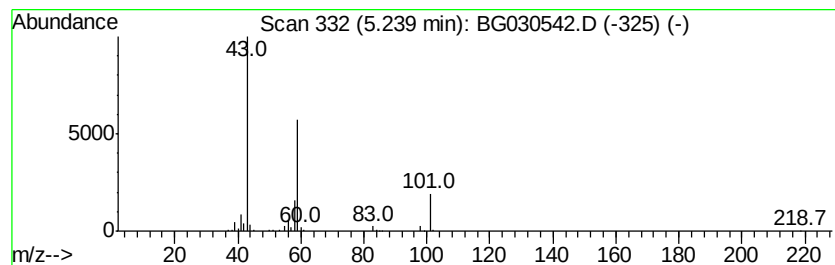
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	15.87 ng	266213	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol	102	C6H14O	000623-37-0	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Propanoic acid, 3-nitro-, methyl...	133	C4H7NO4	020497-95-4	9



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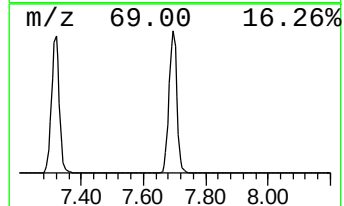
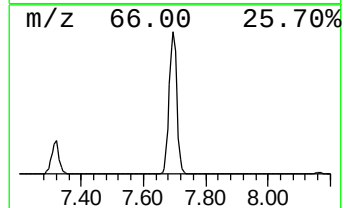
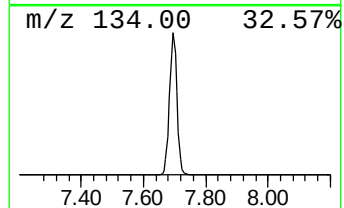
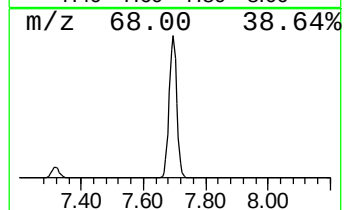
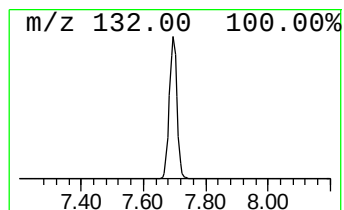
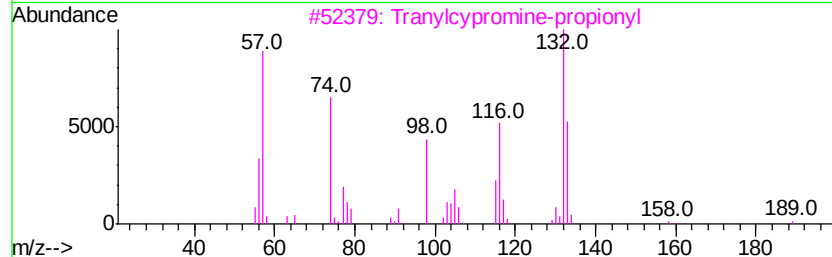
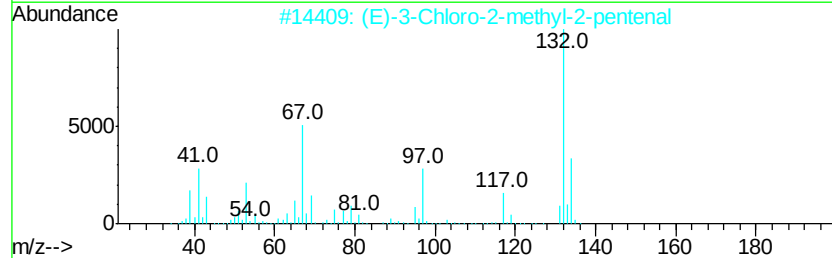
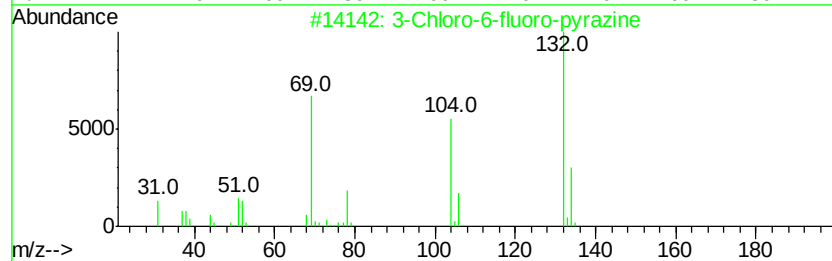
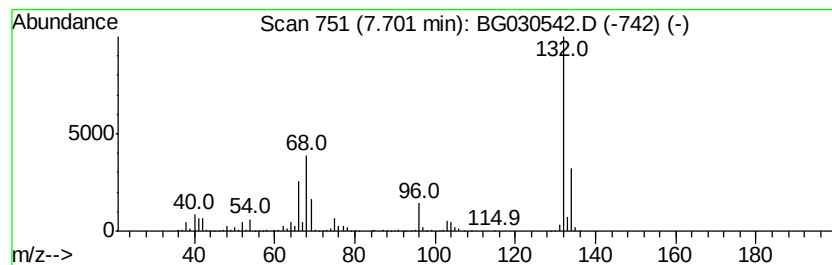
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Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	100.64 ng	1687980	1,4-Dichlorobenzene-d4	8.17

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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	15.9	ng	266213	1	8.17	335442	20.0
unknown7.70	7.70	100.6	ng	1687980	1	8.17	335442	20.0