

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122717\
 Data File : BG031805.D
 Acq On : 27 Dec 2017 20:29
 Operator : SJ/JU
 Sample : I7058-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 122117-SIDI-F-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-BG122117.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	3.434	18	25	36	rBV2	85814	176092	4.53%	1.231%
2	3.522	36	40	51	rVB	39899	65335	1.68%	0.457%
3	6.219	492	499	510	rBV	145271	253296	6.52%	1.770%
4	7.899	776	785	796	rBV	91552	180530	4.65%	1.262%
5	8.311	847	855	865	rBV	305050	557907	14.36%	3.899%
6	8.804	932	939	948	rBB	120649	221498	5.70%	1.548%
7	9.139	988	996	1007	rVB	485115	899530	23.15%	6.287%
8	10.003	1134	1143	1154	rBV	341654	632220	16.27%	4.418%
9	11.689	1421	1430	1440	rBV	180265	331309	8.53%	2.315%
10	14.063	1827	1834	1846	rVB	1430060	2253075	57.99%	15.746%
11	15.432	2061	2067	2076	rVB3	341563	573529	14.76%	4.008%
12	16.906	2311	2318	2330	rBV	781824	1230191	31.66%	8.598%
13	18.164	2521	2532	2544	rBV2	527081	827336	21.29%	5.782%
14	20.696	2957	2963	2973	rBV	2905475	3885350	100.00%	27.154%
15	22.053	3190	3194	3205	rBV3	32954	58809	1.51%	0.411%
16	22.512	3263	3272	3285	rBV	559995	1064748	27.40%	7.441%
17	26.254	3896	3909	3930	rBV2	325772	1097790	28.25%	7.672%

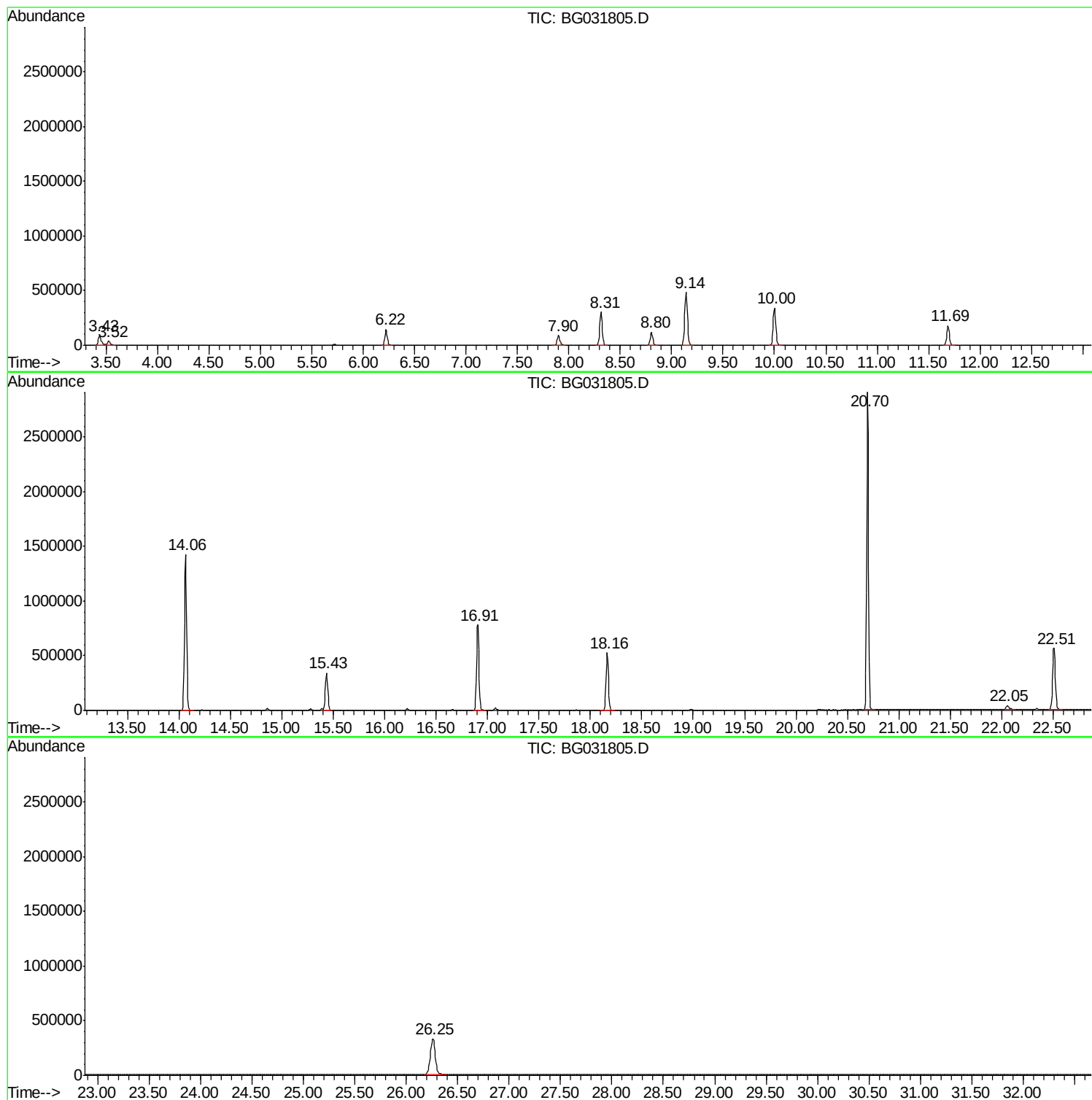
Sum of corrected areas: 14308545

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.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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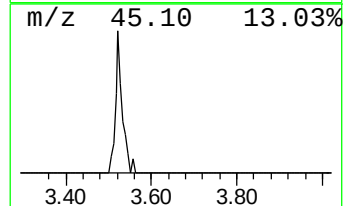
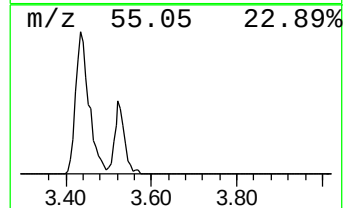
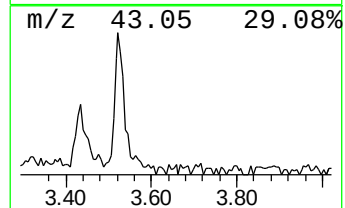
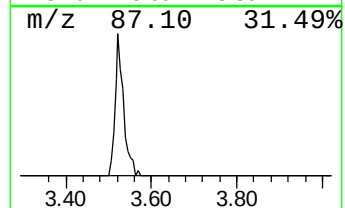
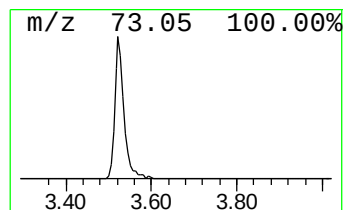
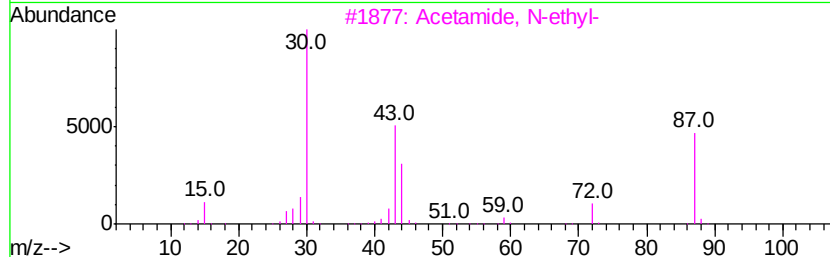
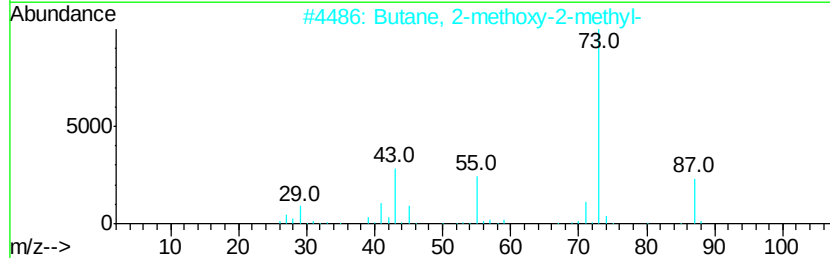
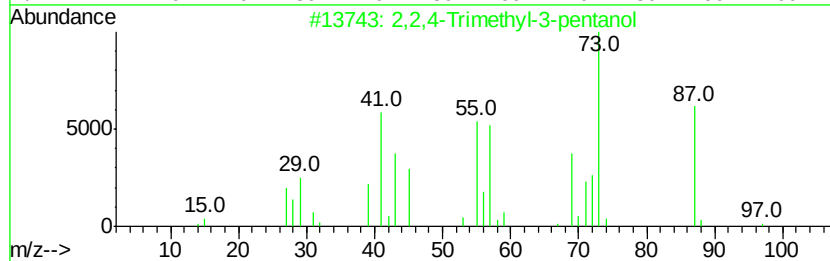
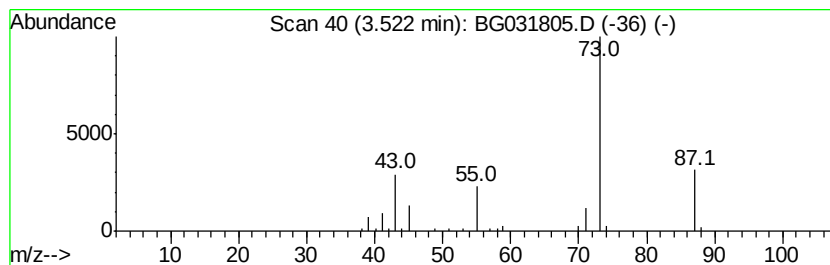
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Peak Number 2 2,2,4-Trimethyl-3-pentanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	5.90 ng	65335	1,4-Dichlorobenzene-d4	8.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	9



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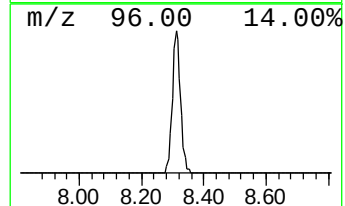
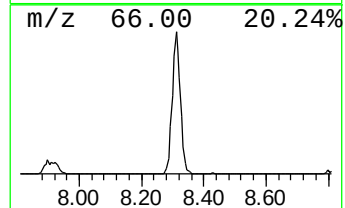
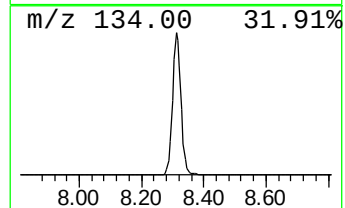
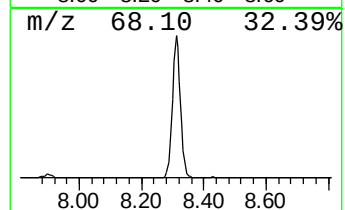
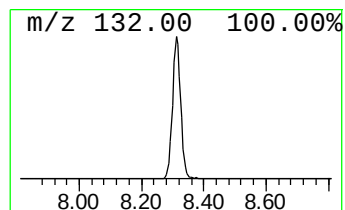
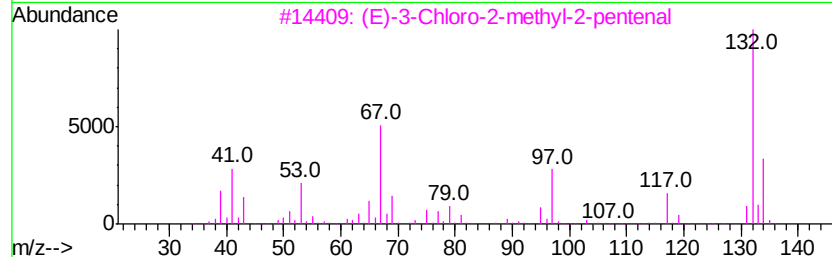
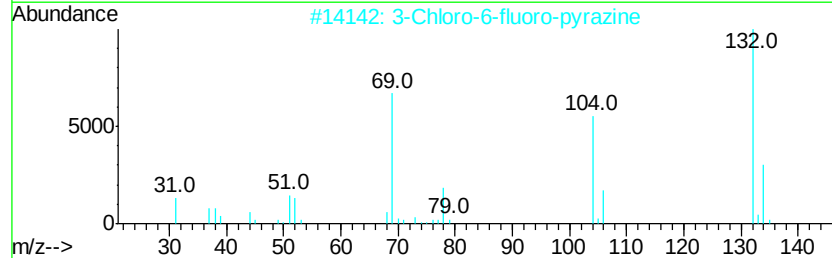
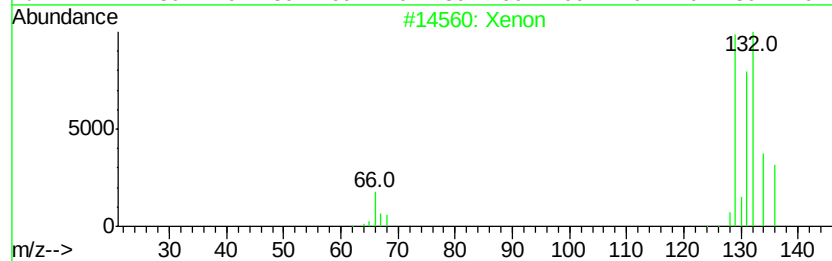
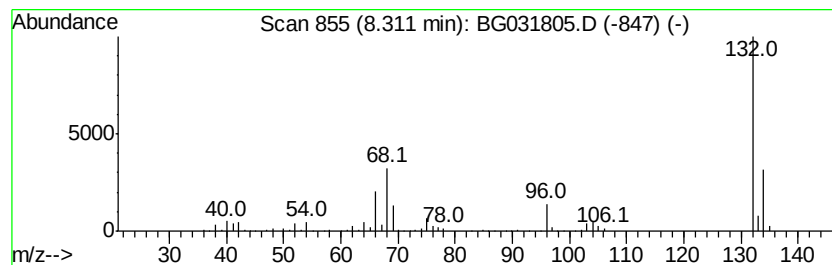
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Peak Number 3 unknown8.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	50.38 ng	557907	1,4-Dichlorobenzene-d4	8.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Xenon	132	Xe	007440-63-3	45
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,2,4-Trimethyl-3...	3.52	5.9	ng	65335	1	8.80	221498	20.0
unknown8.31	8.31	50.4	ng	557907	1	8.80	221498	20.0