

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032819\
Data File : BG040213.D
Acq On : 29 Mar 2019 00:24
Operator : JU/SJ
Sample : PB118375BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB118375BL

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-BG032719.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	3.134	4	8	18	rVB4	27522	50816	1.30%	0.274%
2	5.619	423	431	444	rBV	866243	1411134	36.04%	7.610%
3	7.223	696	704	715	rBV	892342	1529945	39.07%	8.251%
4	7.599	760	768	778	rBV	846726	1460152	37.29%	7.875%
5	8.069	841	848	857	rVB	182487	313678	8.01%	1.692%
6	8.392	894	903	911	rBV	622044	1084904	27.71%	5.851%
7	9.244	1040	1048	1059	rBV	499914	894531	22.84%	4.824%
8	10.884	1318	1327	1335	rBV	244182	432453	11.04%	2.332%
9	13.322	1735	1742	1750	rBV	1348550	2114698	54.00%	11.405%
10	14.697	1970	1976	1985	rBV2	373221	613985	15.68%	3.311%
11	16.183	2223	2229	2252	rBV	1039353	1632369	41.69%	8.803%
12	17.446	2437	2444	2452	rBV2	566334	860127	21.97%	4.639%
13	20.043	2880	2886	2892	rBV	3030787	3915748	100.00%	21.118%
14	21.724	3164	3172	3181	rBV	679650	1135641	29.00%	6.125%
15	25.020	3722	3733	3745	rVB2	376339	1092333	27.90%	5.891%

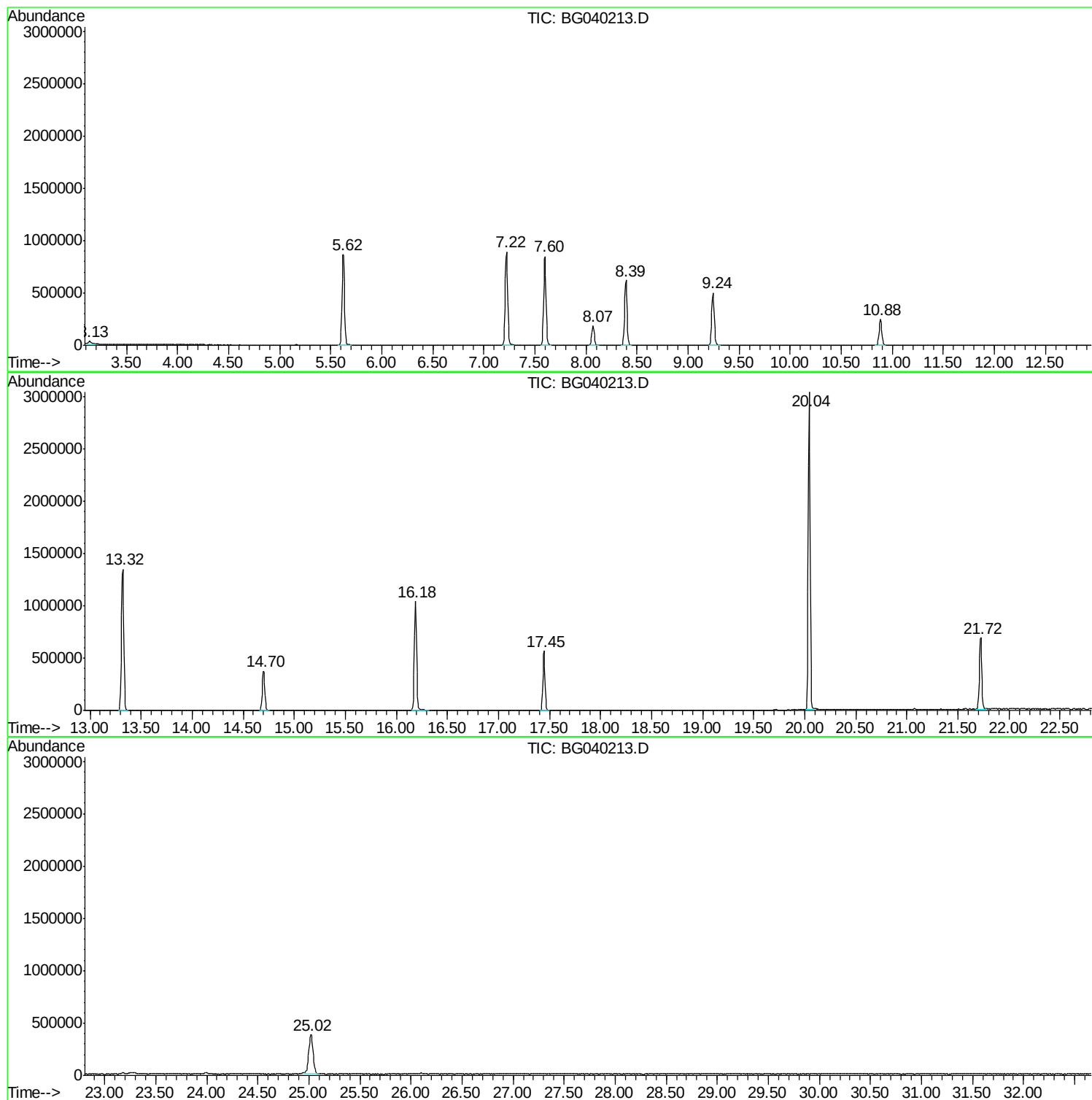
Sum of corrected areas: 18542514

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.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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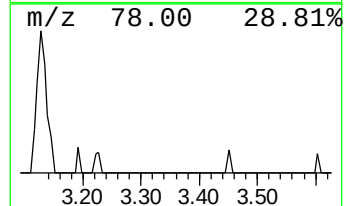
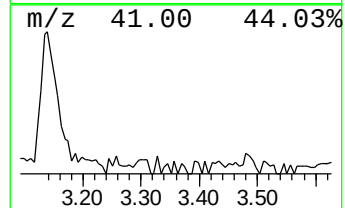
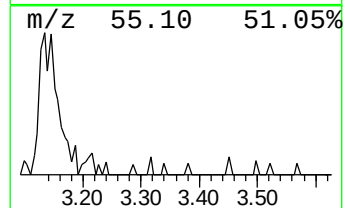
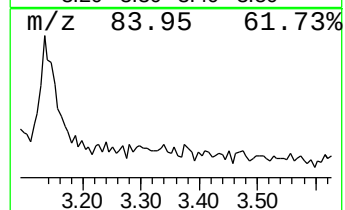
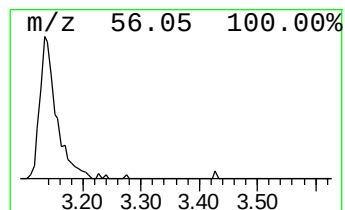
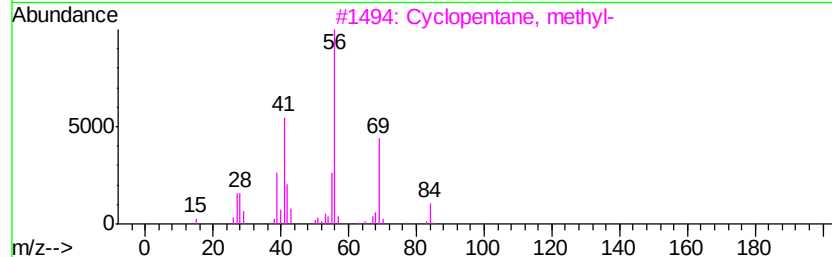
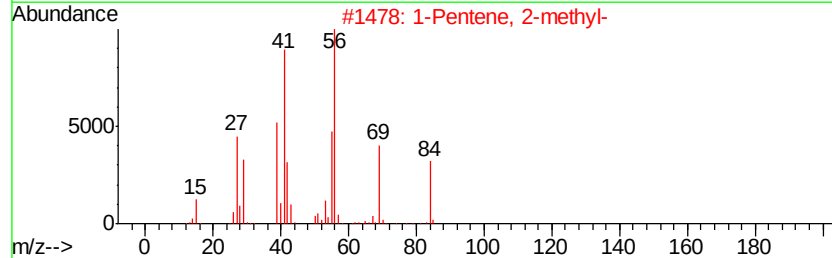
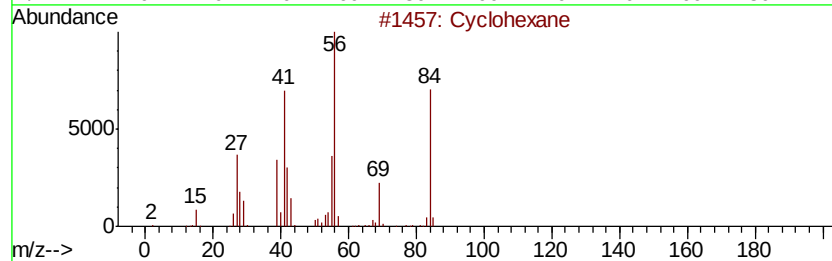
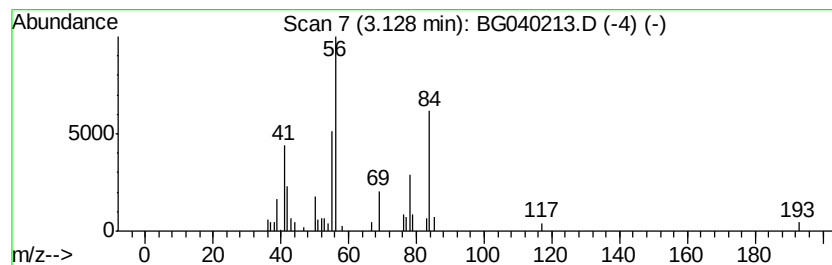
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Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	3.24 ng	50816	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	87
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	56
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	42
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	37
5		2-Hexene, 5-methyl-, (E)-	98	C7H14	007385-82-2	37



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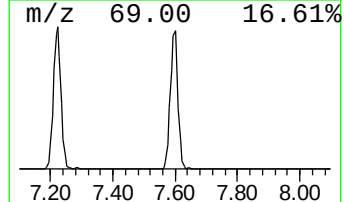
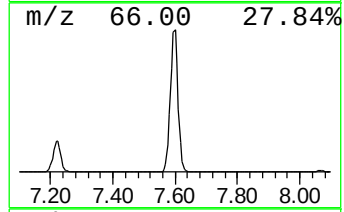
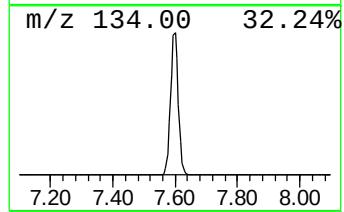
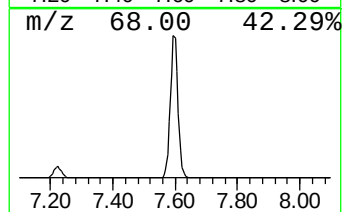
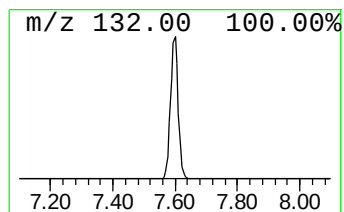
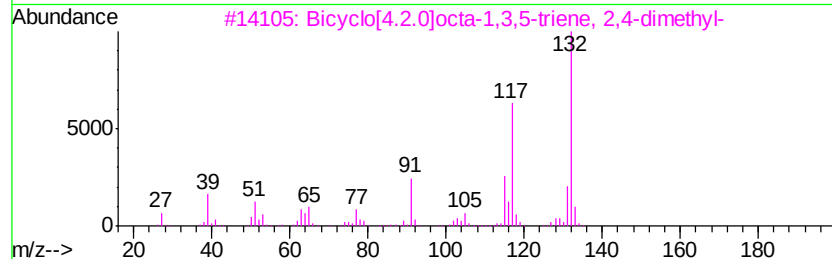
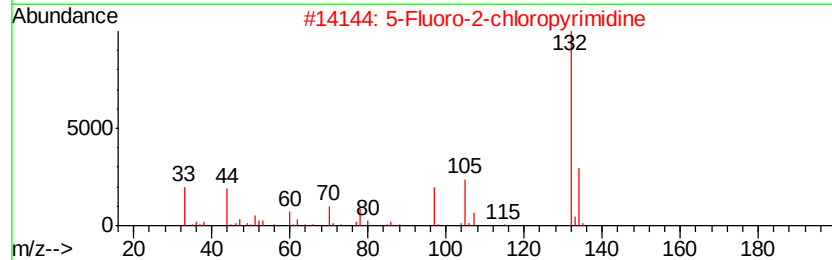
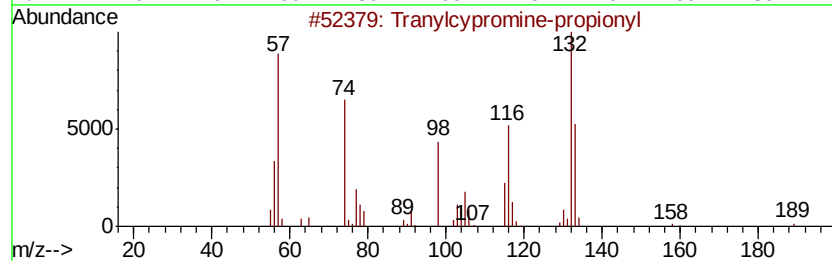
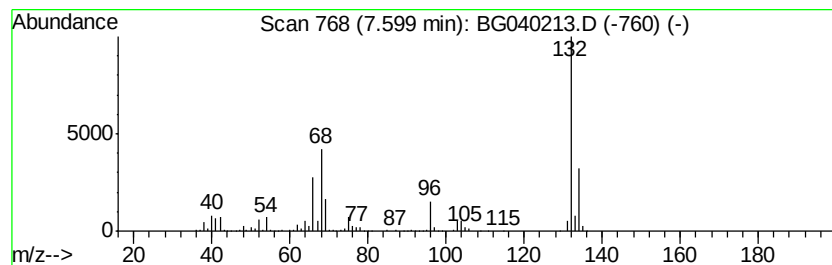
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	93.10 ng	1460150	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	3.13	3.2	ng	50816	1	8.07	313678	20.0
unknown7.60	7.60	93.1	ng	1460150	1	8.07	313678	20.0