

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040119\
 Data File : BG040270.D
 Acq On : 1 Apr 2019 18:18
 Operator : JU/SJ
 Sample : PB118456BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118456BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	3.122	3	6	15	rVB	26331	44060	1.16%	0.226%
2	5.608	422	429	442	rBV	841426	1312435	34.53%	6.740%
3	7.206	693	701	712	rBV	859156	1451183	38.18%	7.453%
4	7.582	756	765	778	rBV	821508	1388469	36.53%	7.131%
5	8.052	837	845	854	rBV	196660	335088	8.82%	1.721%
6	8.375	893	900	908	rBV	568815	1003776	26.41%	5.155%
7	9.227	1035	1045	1059	rBV	472935	844560	22.22%	4.337%
8	10.866	1317	1324	1335	rBV	269098	487780	12.83%	2.505%
9	13.305	1732	1739	1747	rBV	1445041	2191202	57.65%	11.253%
10	14.685	1967	1974	1985	rVB2	493450	788784	20.75%	4.051%
11	16.172	2220	2227	2241	rBV	1112464	1791573	47.14%	9.201%
12	17.429	2434	2441	2450	rBV	746458	1134436	29.85%	5.826%
13	20.032	2878	2884	2890	rBV	2834774	3800924	100.00%	19.520%
14	21.706	3161	3169	3177	rBV	782662	1298797	34.17%	6.670%
15	22.470	3293	3299	3304	rBV2	37256	58676	1.54%	0.301%
16	23.169	3412	3418	3424	rBV2	42032	78848	2.07%	0.405%
17	23.986	3550	3557	3565	rVB2	39528	86765	2.28%	0.446%
18	25.003	3715	3730	3742	rVB	428159	1319144	34.71%	6.775%
19	26.101	3911	3917	3928	rVB4	17975	55123	1.45%	0.283%

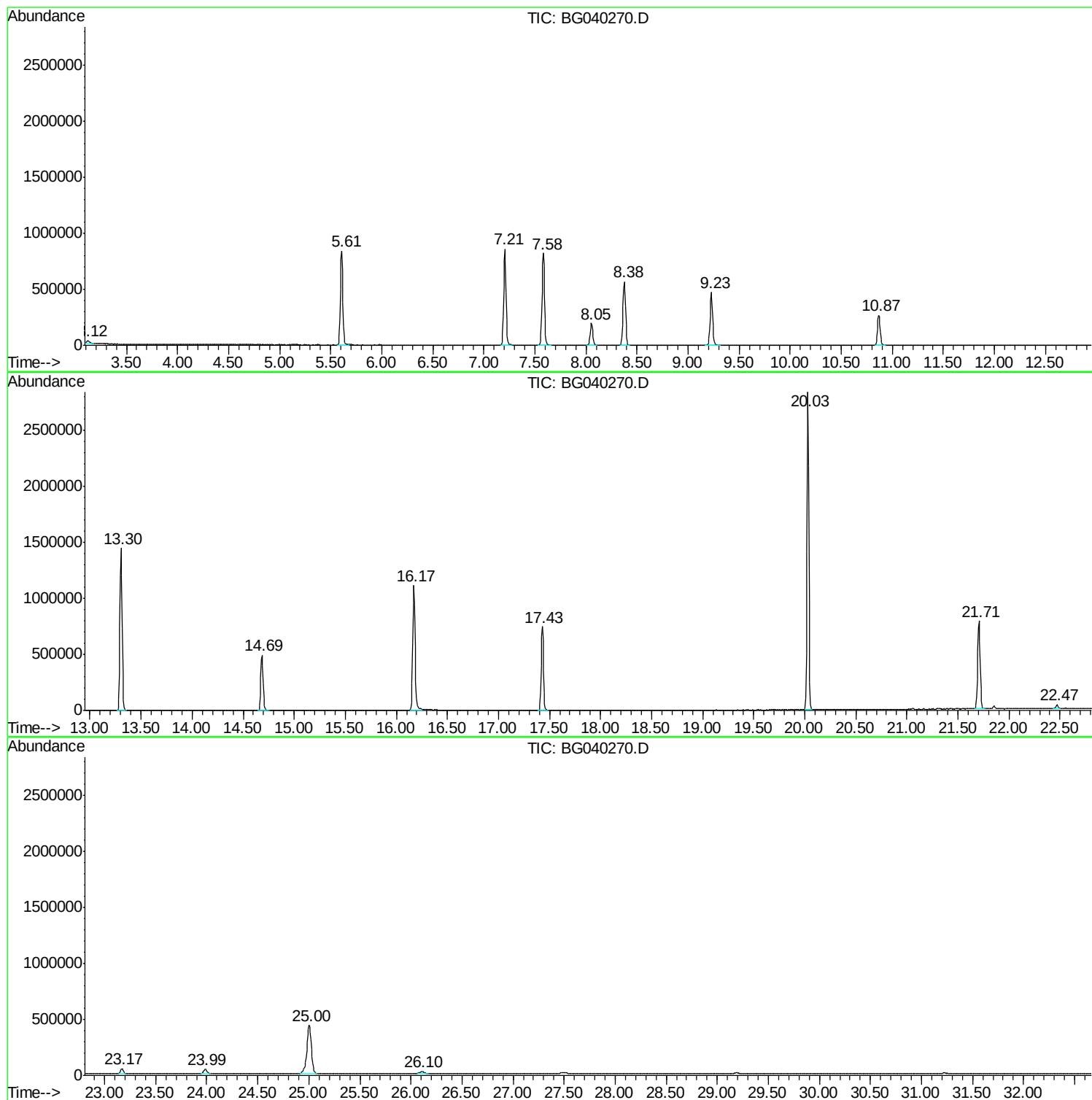
Sum of corrected areas: 19471623

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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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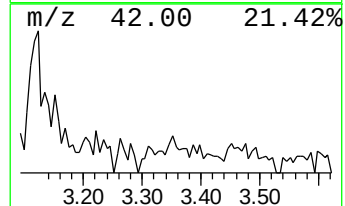
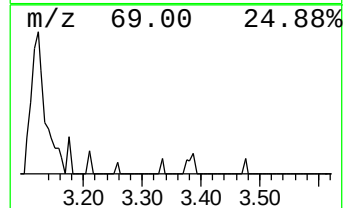
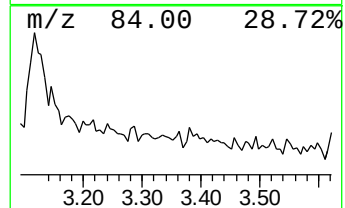
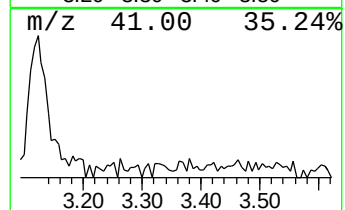
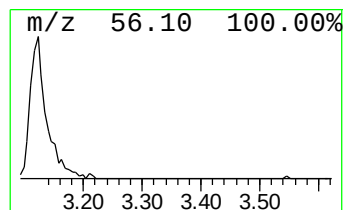
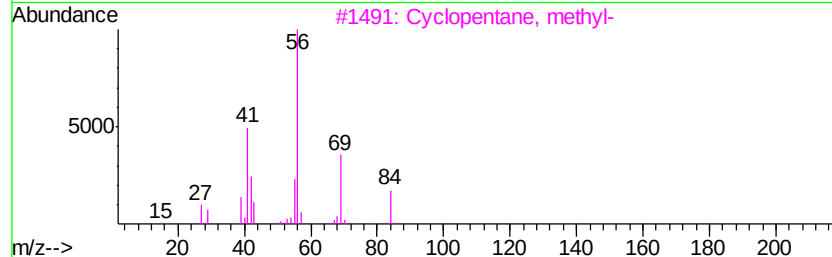
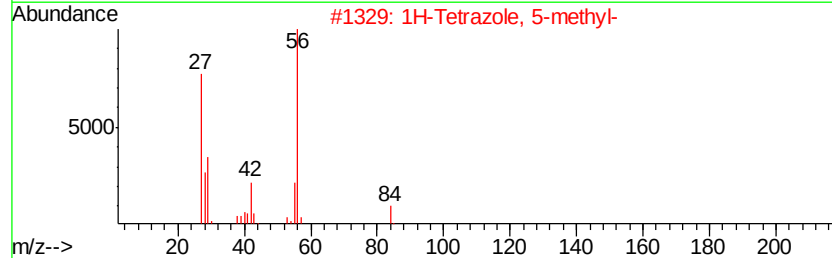
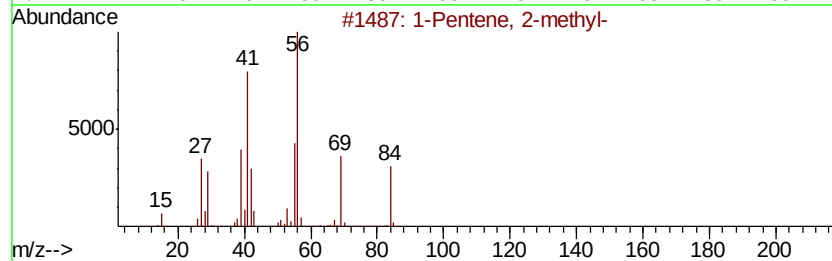
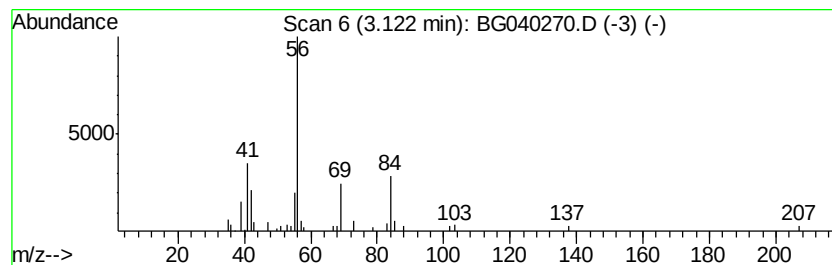
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Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.12	2.63 ng	44060	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	53
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	53
4		Cyclohexane	84	C6H12	000110-82-7	50
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	47



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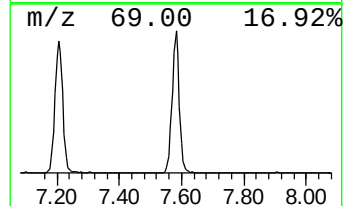
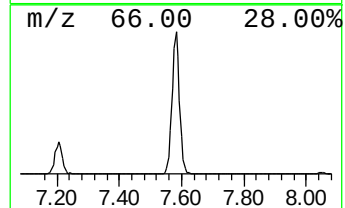
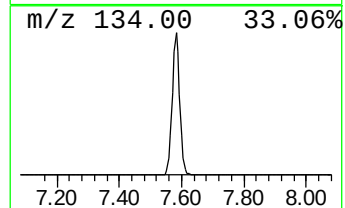
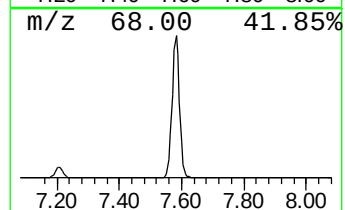
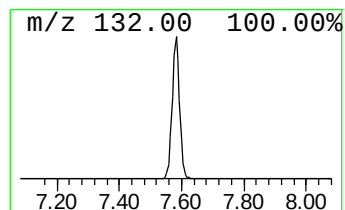
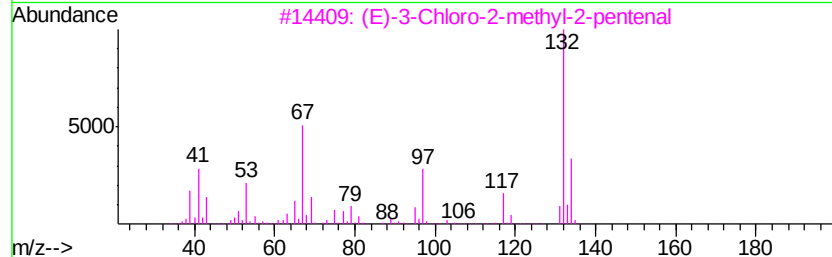
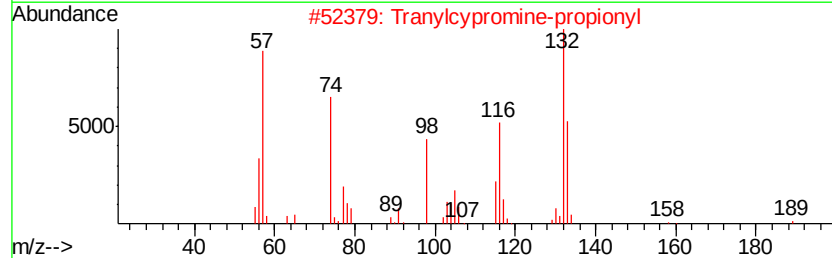
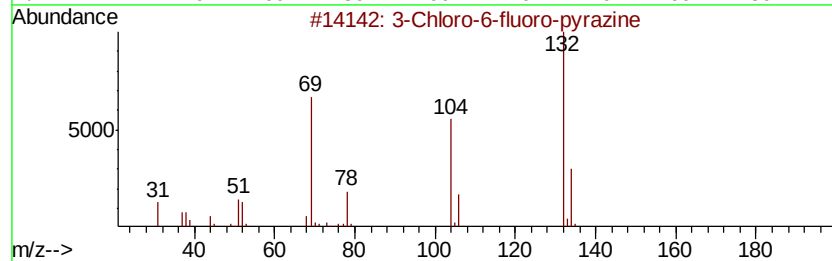
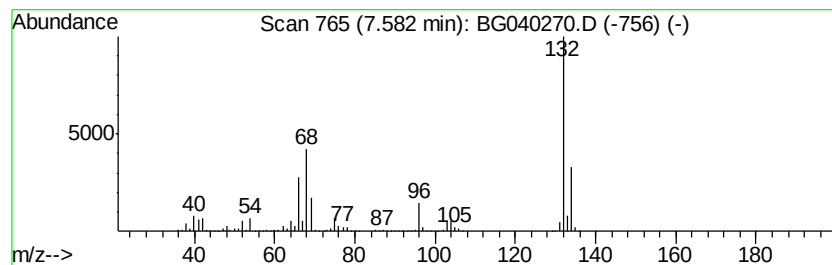
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	82.87 ng	1388470	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	3.12	2.6	ng	44060	1	8.05	335088	20.0
unknown7.58	7.58	82.9	ng	1388470	1	8.05	335088	20.0