

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031471.D
 Acq On : 12 Dec 2017 1:57
 Operator : SJ/JU
 Sample : I6762-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120417-TIDO-F-U-B

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-BG121117.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	5.187	316	323	334	rBV	45174	78691	1.46%	0.436%
2	5.669	399	405	418	rBV	237947	444650	8.28%	2.462%
3	7.279	672	679	693	rBV	143940	303917	5.66%	1.683%
4	7.643	734	741	760	rBV	459838	849576	15.81%	4.703%
5	8.107	813	820	830	rBV	165757	295639	5.50%	1.637%
6	8.430	868	875	889	rBV	640290	1180041	21.96%	6.533%
7	9.276	1012	1019	1034	rBV	444448	893748	16.63%	4.948%
8	10.927	1291	1300	1320	rBV	178798	419486	7.81%	2.322%
9	13.354	1706	1713	1732	rBV	1558645	2585885	48.13%	14.316%
10	14.735	1940	1948	1961	rBV2	371095	675218	12.57%	3.738%
11	16.215	2194	2200	2219	rBV	767182	1362847	25.37%	7.545%
12	17.473	2408	2414	2431	rBV2	595332	986590	18.36%	5.462%
13	20.064	2849	2855	2865	rBV	3931216	5372851	100.00%	29.745%
14	21.356	3070	3075	3088	rBV8	23583	79937	1.49%	0.443%
15	21.744	3134	3141	3156	rBV2	745851	1301567	24.22%	7.206%
16	25.023	3685	3699	3713	rBV2	397769	1232327	22.94%	6.822%

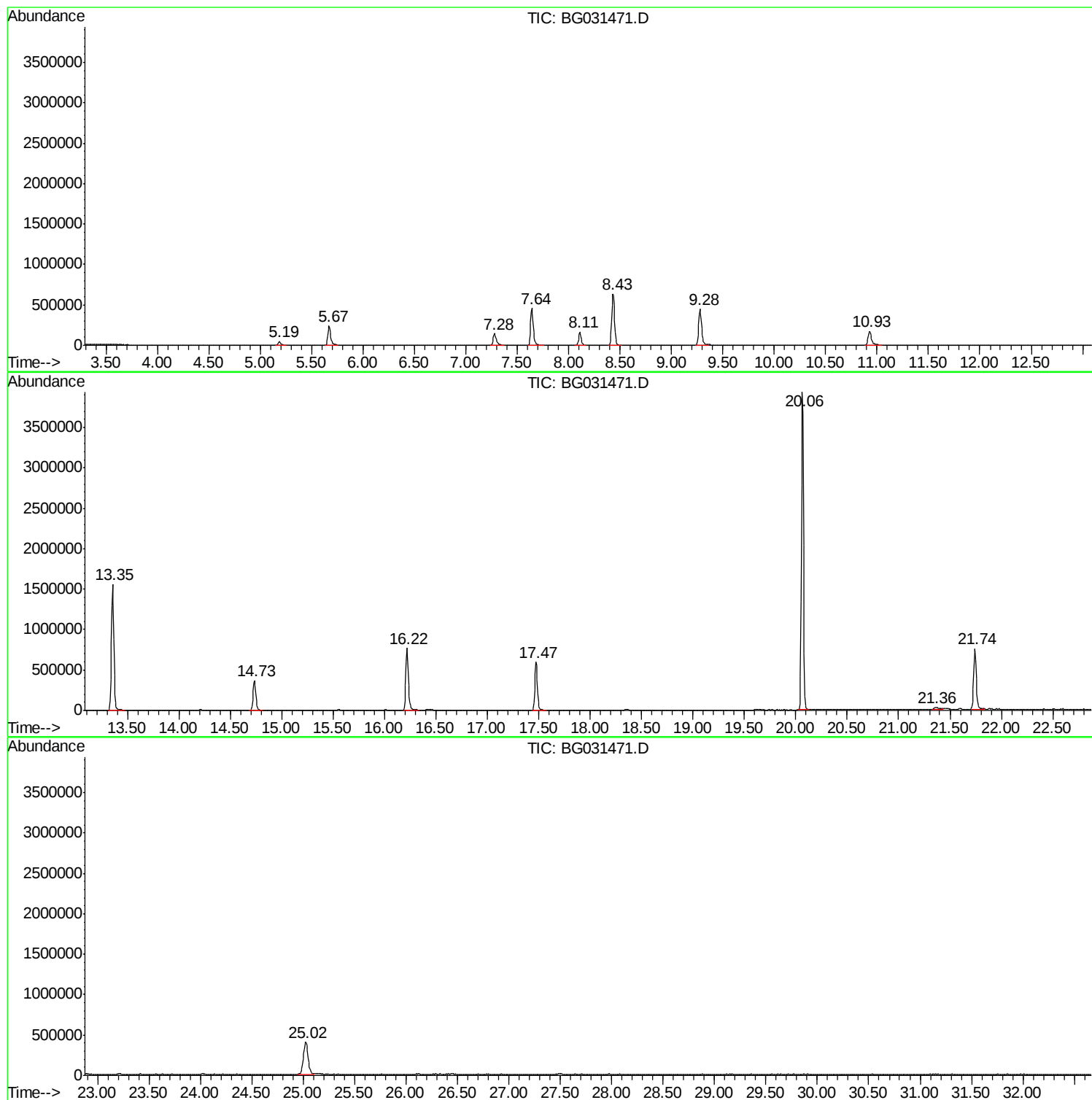
Sum of corrected areas: 18062970

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.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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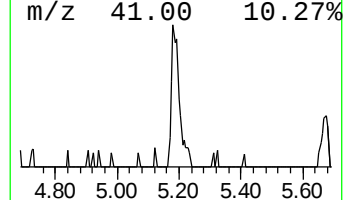
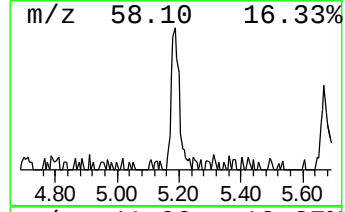
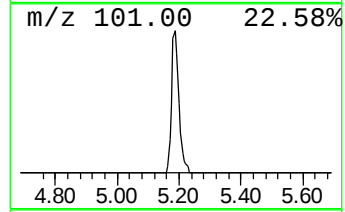
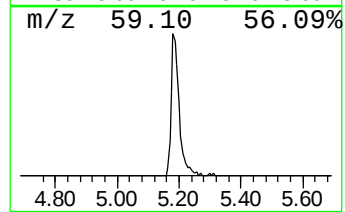
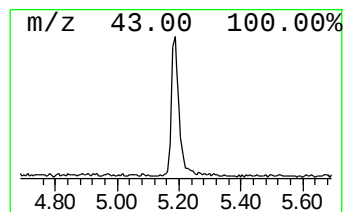
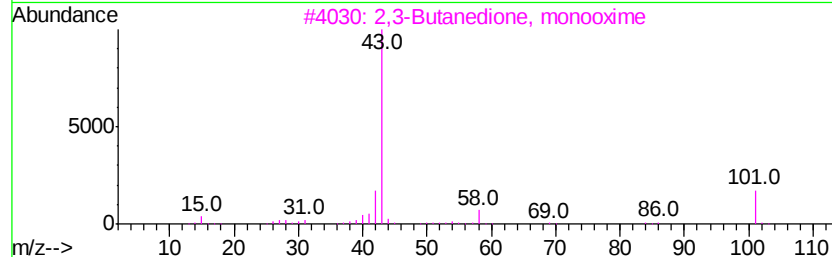
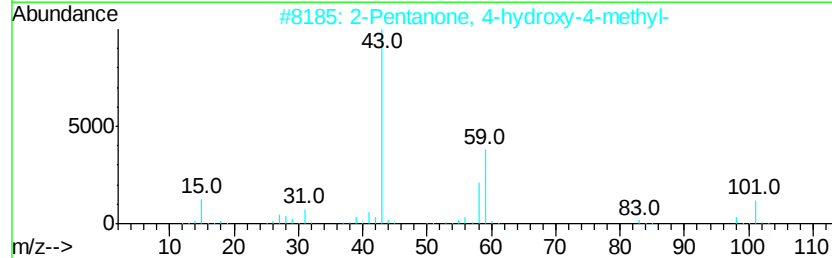
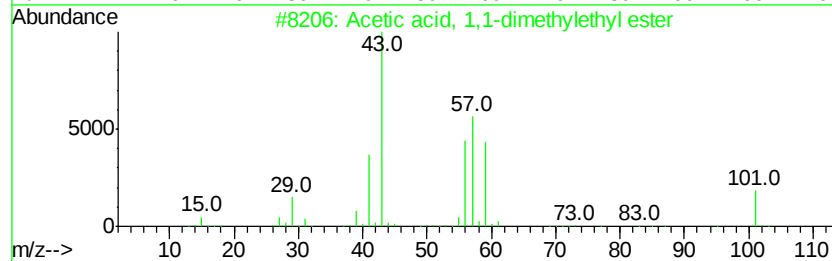
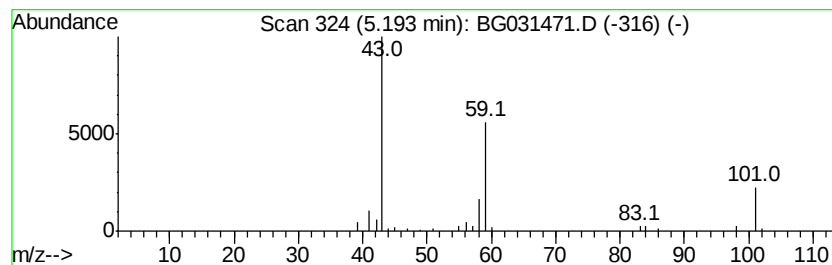
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Acetic acid, 1,1-dimethylet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	5.32 ng	78691	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	17
5		Propanoic acid, 2-hydroxy-2-meth...	132	C6H12O3	000080-55-7	17



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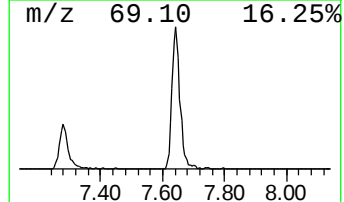
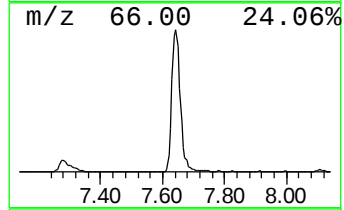
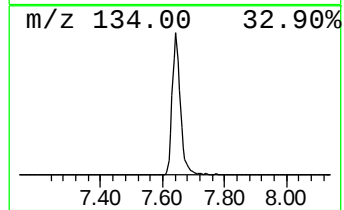
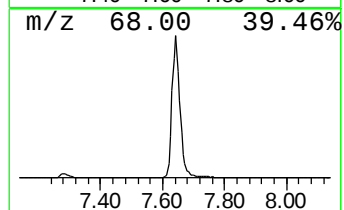
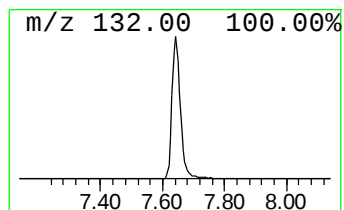
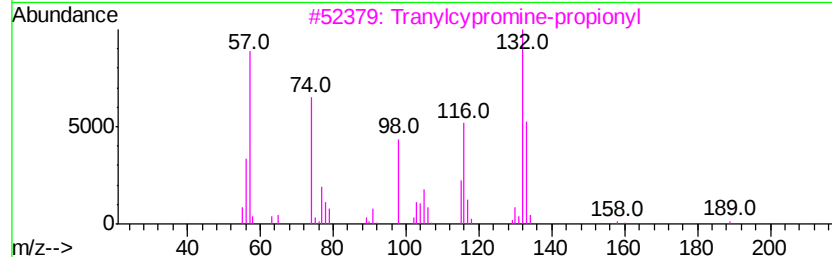
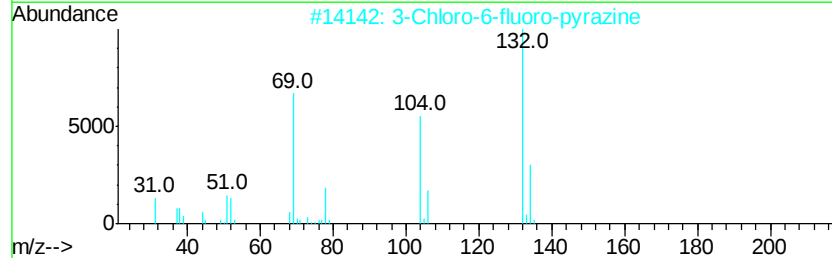
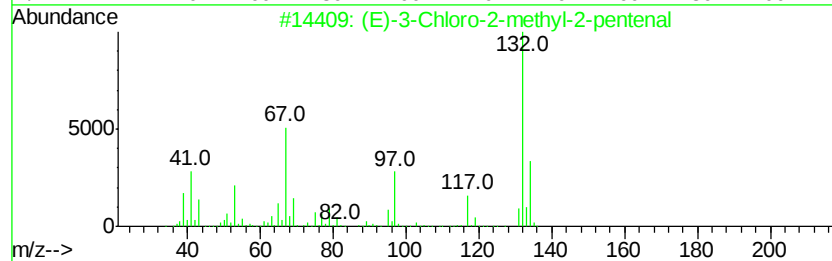
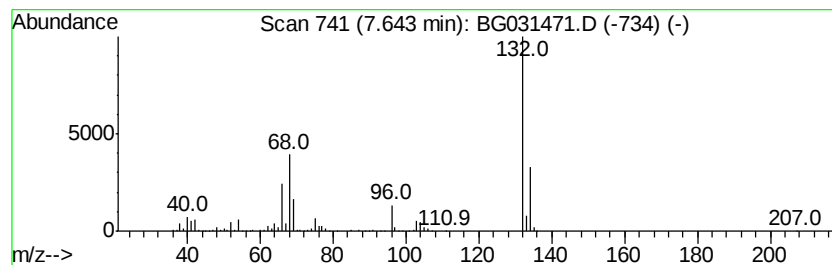
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	57.47 ng	849576	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid, 1,1-...	5.19	5.3	ng	78691	1	8.11	295639	20.0
unknown7.64	7.64	57.5	ng	849576	1	8.11	295639	20.0