

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF013118\
 Data File : BF102683.D
 Acq On : 1 Feb 2018 4:11
 Operator : SJ/JU
 Sample : PB106224BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB106224BL

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_F\METHODS\8270-BF012218.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	4.916	477	481	492	rBV	167919	247572	7.32%	0.966%
2	5.339	549	553	573	rBV	3103403	3382338	100.00%	13.200%
3	6.369	724	728	744	rBV	2955694	3240826	95.82%	12.648%
4	6.486	744	748	763	rVB	3437954	3300135	97.57%	12.880%
5	6.704	782	785	795	rVB	880106	771298	22.80%	3.010%
6	6.863	808	812	815	rBV	2701631	2413997	71.37%	9.421%
7	7.275	878	882	899	rBV	1667589	1813780	53.63%	7.079%
8	7.986	999	1003	1015	rBV	853381	930550	27.51%	3.632%
9	9.063	1181	1186	1200	rBV	3183791	2903107	85.83%	11.330%
10	9.739	1297	1301	1310	rBV	810020	804687	23.79%	3.140%
11	10.533	1432	1436	1452	rBV	1239971	1335348	39.48%	5.212%
12	11.227	1550	1554	1569	rBV	756041	732062	21.64%	2.857%
13	12.621	1787	1791	1794	rBV	57055	50194	1.48%	0.196%
14	12.810	1819	1823	1826	rBV	2708718	2377388	70.29%	9.278%
15	13.868	1999	2003	2010	rBV	795801	787593	23.29%	3.074%
16	15.298	2241	2246	2253	rBV	377122	532165	15.73%	2.077%

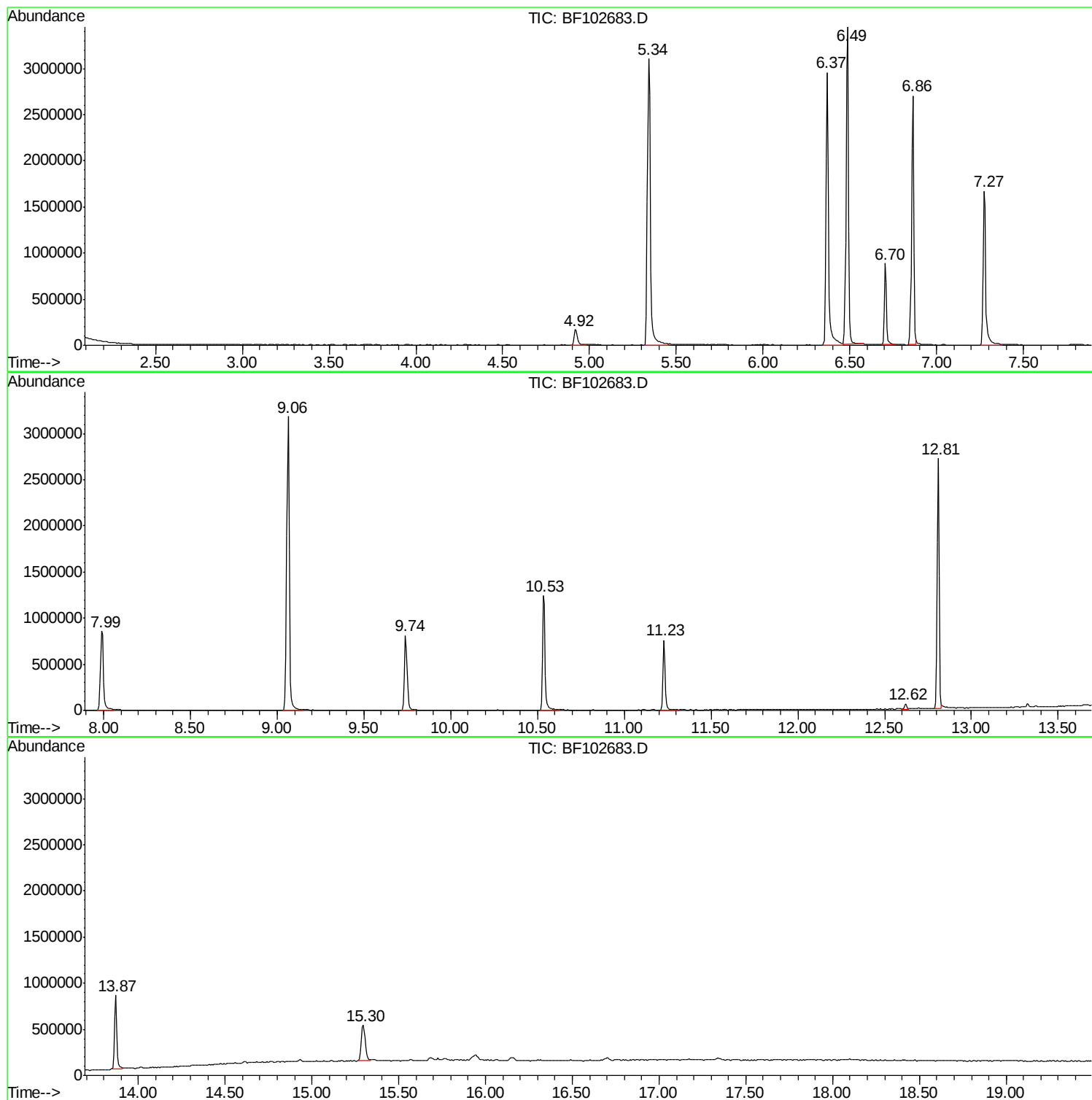
Sum of corrected areas: 25623040

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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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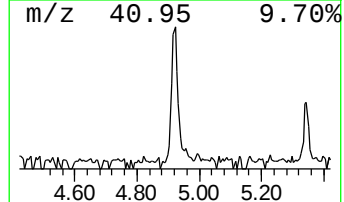
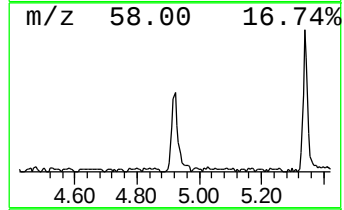
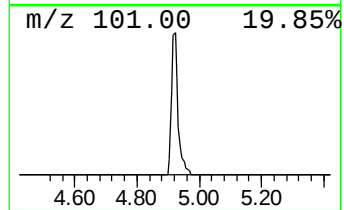
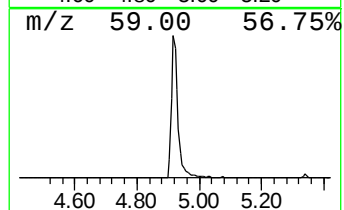
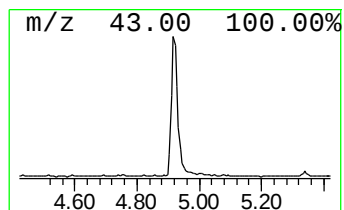
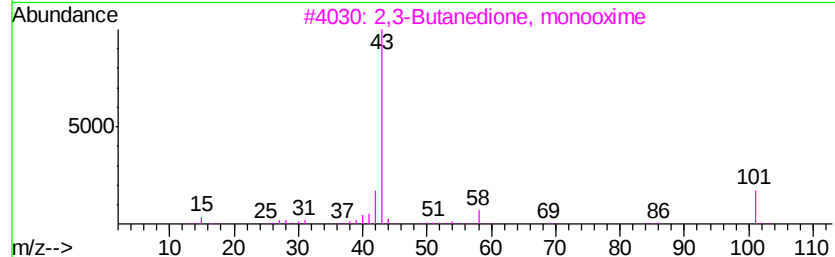
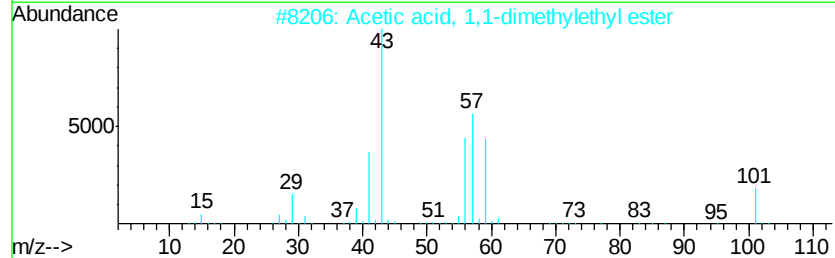
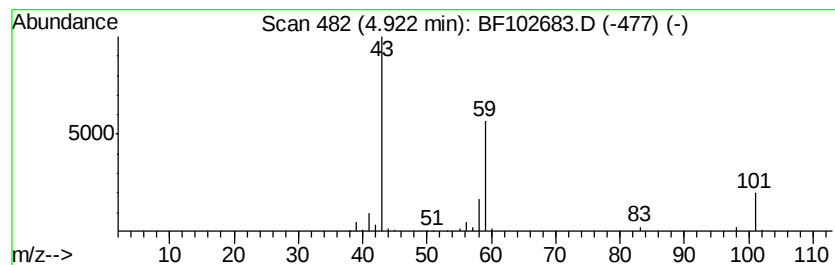
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	6.42 ng	247572	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			1,2-Butanediol	90	C4H10O2	000584-03-2	9



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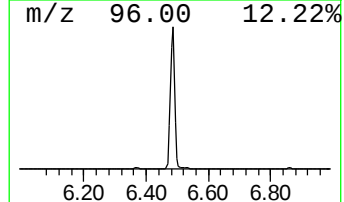
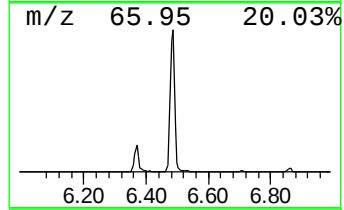
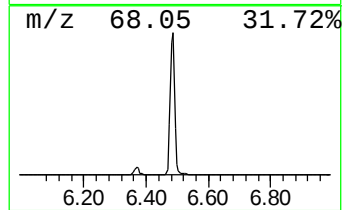
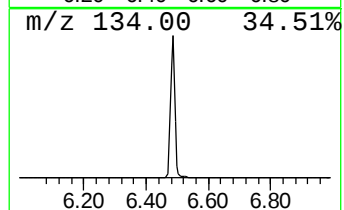
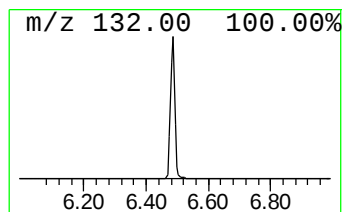
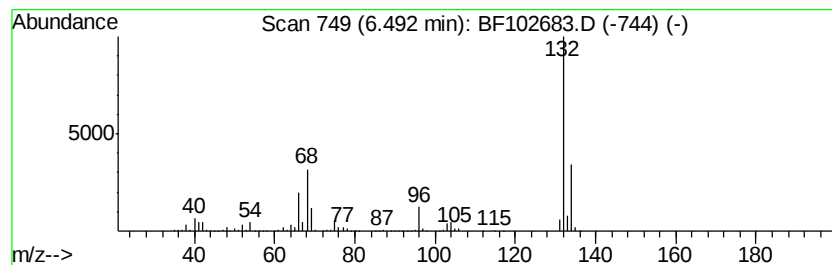
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Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	85.57 ng	3300140	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.92	6.4	ng	247572	1	6.70	771298	20.0
unknown6.49	6.49	85.6	ng	3300140	1	6.70	771298	20.0