

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122117\
 Data File : BF101631.D
 Acq On : 22 Dec 2017 3:34
 Operator : SJ/JU
 Sample : I7008-01
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 1 % 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-BF121417.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	5.004	493	496	512	rBV	133912	210140	6.50%	1.037%
2	5.416	563	566	591	rBV	838056	1260399	39.01%	6.219%
3	6.439	737	740	758	rBV	532917	866720	26.82%	4.277%
4	6.569	758	762	769	rBV	1901467	1987043	61.49%	9.805%
5	6.792	797	800	810	rBV	826863	804859	24.91%	3.971%
6	6.951	823	827	844	rBV	2624349	2540478	78.62%	12.536%
7	7.369	893	898	922	rBV	1863222	2182213	67.53%	10.768%
8	8.075	1015	1018	1036	rBV	916157	1035449	32.04%	5.109%
9	9.151	1197	1201	1204	rBV	3447946	3231350	100.00%	15.945%
10	9.833	1314	1317	1329	rVB	901794	911142	28.20%	4.496%
11	10.627	1448	1452	1465	rBV	935293	1018879	31.53%	5.028%
12	11.321	1567	1570	1581	rBV	727423	723162	22.38%	3.568%
13	12.904	1835	1839	1842	rBV	2342515	2141144	66.26%	10.565%
14	13.792	1986	1990	2001	rBV5	24031	53526	1.66%	0.264%
15	13.968	2016	2020	2030	rBV	730603	745116	23.06%	3.677%
16	15.433	2265	2269	2279	rBV	391008	554274	17.15%	2.735%

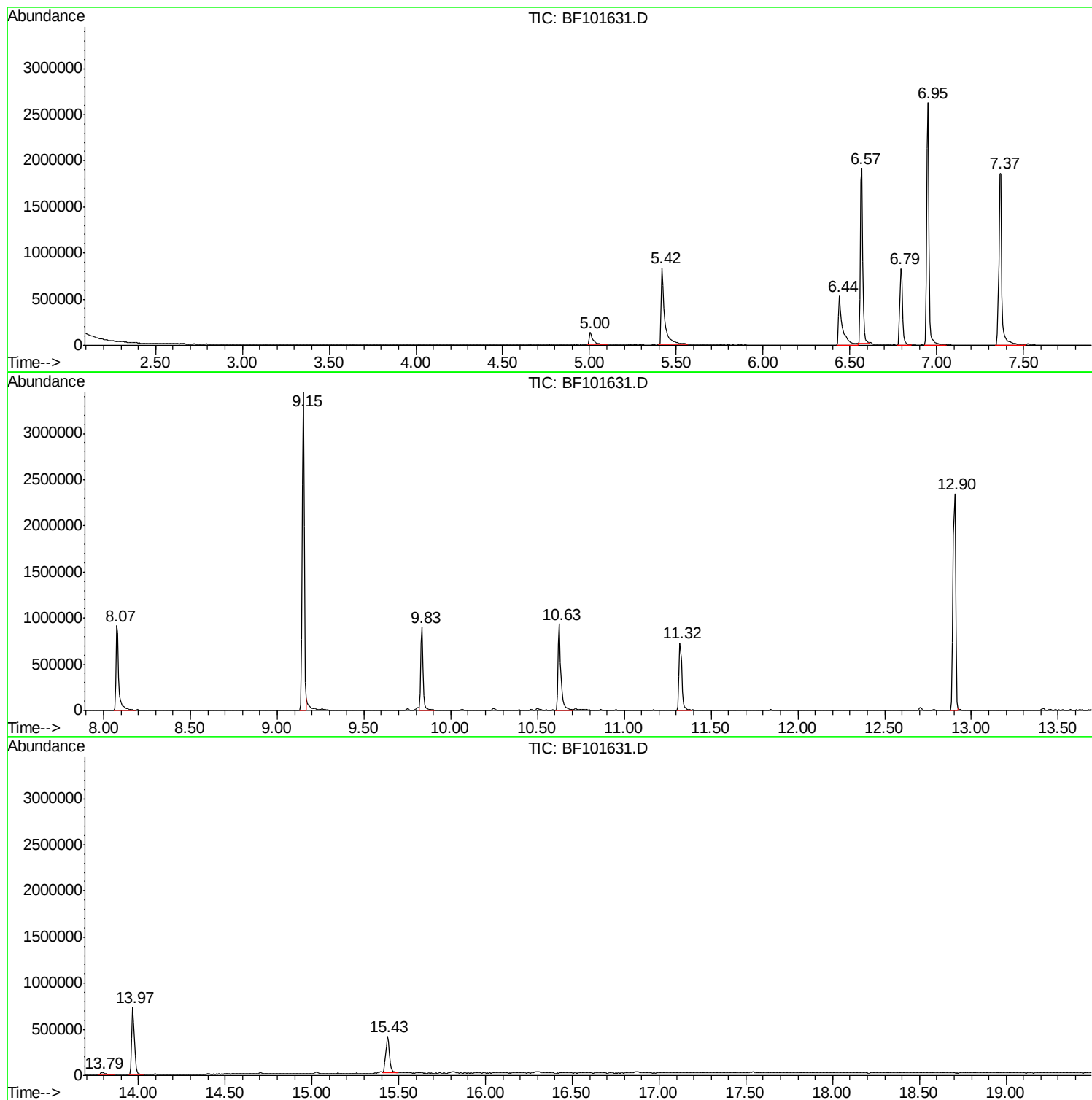
Sum of corrected areas: 20265894

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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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.
5.00	5.22 ng	210140	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		

 Peak Number 2 3-Chloro-6-fluoro-pyrazine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	49.38 ng	1987040	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	(E)-3	Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16	
3	5	Aminoindole	132	C8H8N2	005192-03-0	12	
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10		
5	1,3,5	Trifluorobenzene	132	C6H3F3	000372-38-3	9	

 Peak Number 3 1,4-Dichlorobenzene-D4 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.95	63.13 ng	2540480	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dichlorobenzene-D4	150	C6D4Cl2	003855-82-1	97
2			1,2-Dichlorobenzene-D4	150	C6D4Cl2	002199-69-1	94
3			9-Borabicyclo[3.3.1]nonane, 9-et...	150	C10H19B	052102-17-7	38
4			2,5-Cyclohexadiene-1,4-dione, 3-...	362	C17H15ClN2O5	313480-08-9	14
5			2-Bromo-2'-nitroacetophenone	243	C8H6BrN03	006851-99-6	12

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	5.2	ng	210140	1	6.79	804859	20.0
3-Chloro-6-fluoro...	6.57	49.4	ng	1987040	1	6.79	804859	20.0
1,4-Dichlorobenze...	6.95	63.1	ng	2540480	1	6.79	804859	20.0