

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100719\
 Data File : BP000539.D
 Acq On : 07 Oct 2019 13:41
 Operator : JU
 Sample : PB123680BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB123680BL

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_P\METHODS\8270-BP100119.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	5.387	557	561	564	rBV	4892531	4505728	52.16%	8.456%
2	6.399	729	733	736	rBV	5065980	4668246	54.04%	8.761%
3	6.534	752	756	759	rBV	5719225	4973135	57.57%	9.333%
4	6.757	791	794	798	rBV	1616737	1324851	15.34%	2.486%
5	6.916	817	821	824	rBV	5413969	4317177	49.97%	8.102%
6	7.316	885	889	892	rBV	3280948	2943835	34.08%	5.525%
7	8.040	1009	1012	1016	rBV	1918323	1640967	19.00%	3.080%
8	9.122	1192	1196	1199	rBV	8122910	6804493	78.77%	12.770%
9	9.804	1308	1312	1316	rBV	2636639	2069906	23.96%	3.885%
10	10.598	1443	1447	1475	rBV	4960871	4387050	50.78%	8.233%
11	11.298	1563	1566	1570	rBV	2526756	2236316	25.89%	4.197%
12	12.910	1835	1840	1843	rBV	8901300	8638856	100.00%	16.213%
13	13.969	2016	2020	2033	rBV	2757825	2370672	27.44%	4.449%
14	15.445	2265	2271	2277	rBV	1924089	2403919	27.83%	4.511%

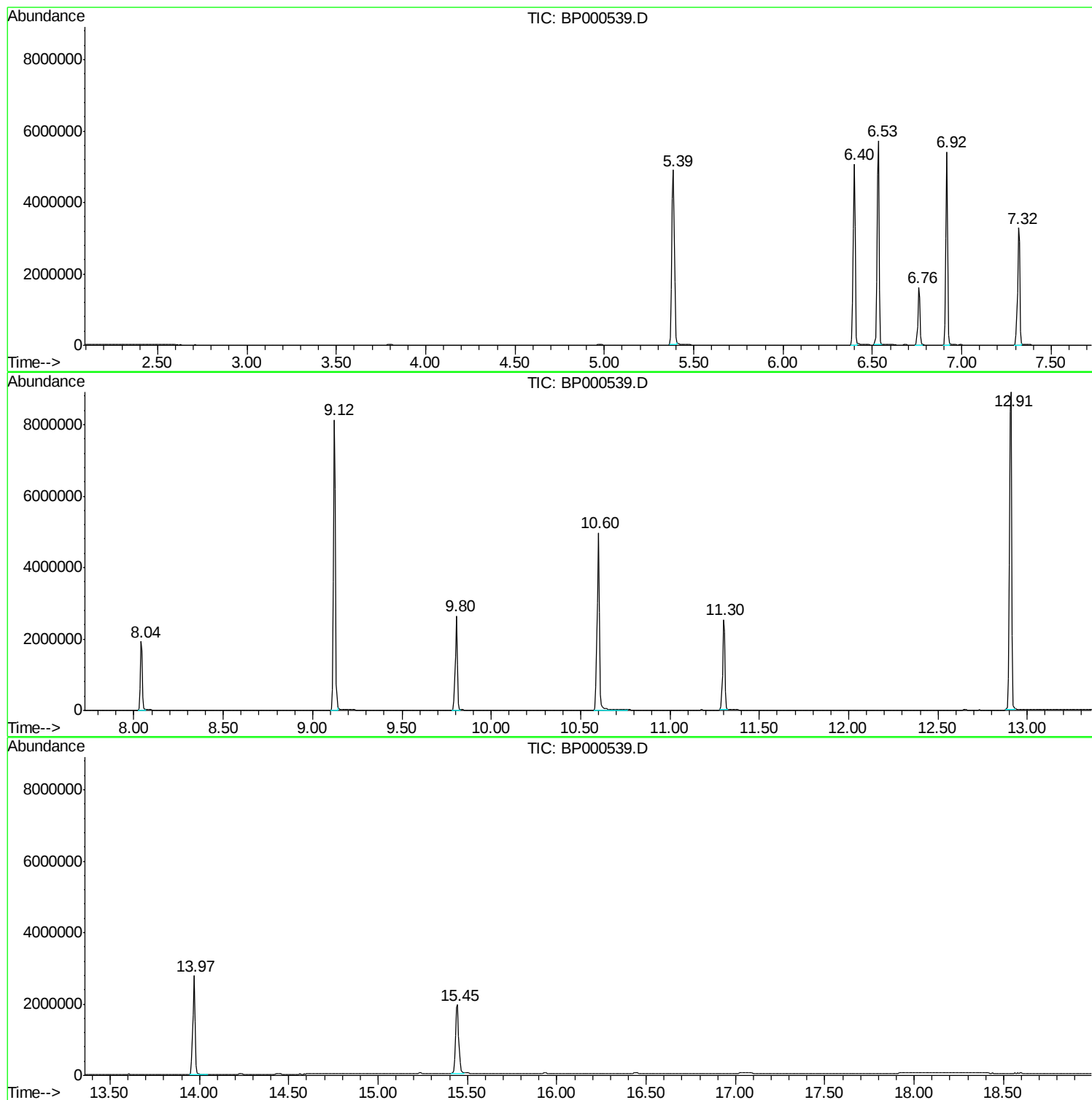
Sum of corrected areas: 53285151

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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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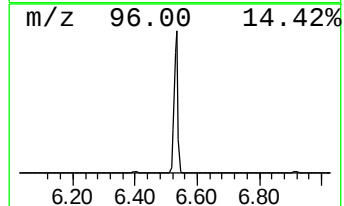
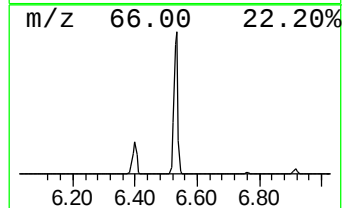
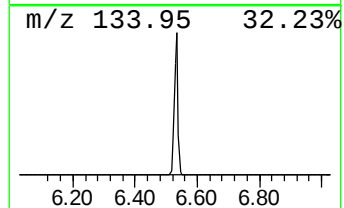
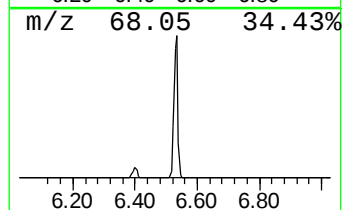
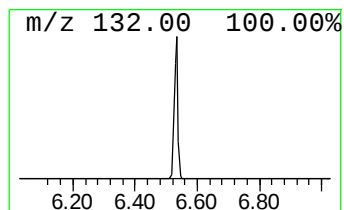
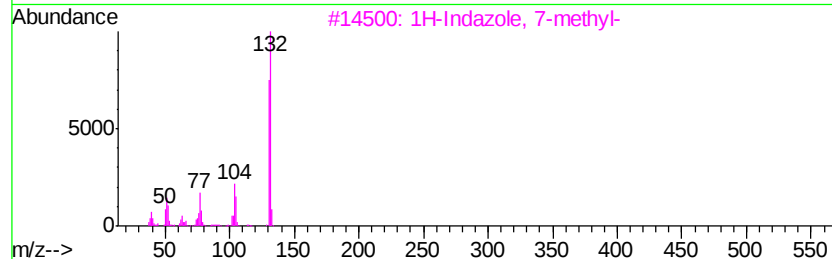
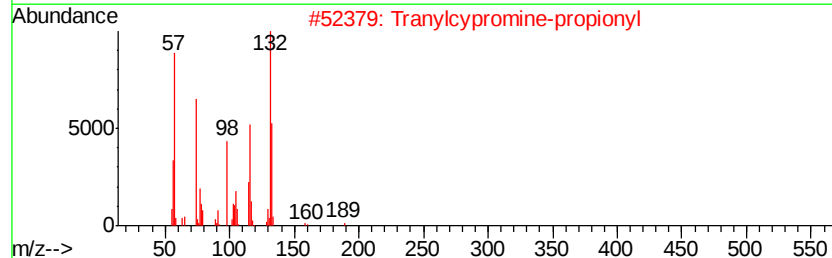
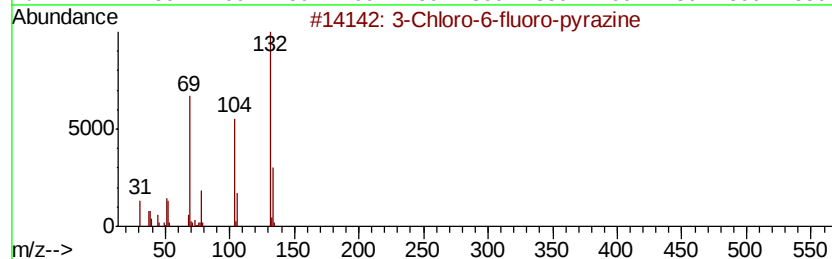
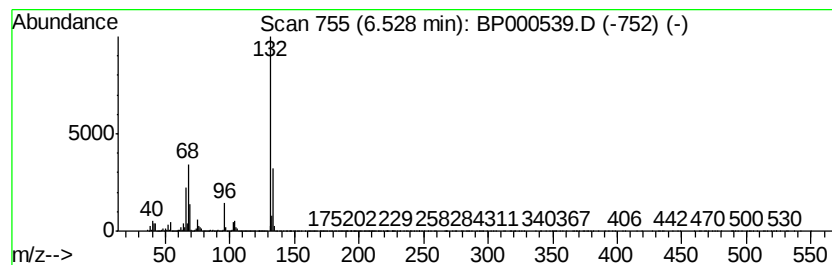
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	75.07 ng	4973140	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	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
unknown6.53	6.53	75.1	ng	4973140	1	6.76	1324850	20.0