

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092019\
Data File : BP000345.D
Acq On : 20 Sep 2019 10:16
Operator : HP/JU
Sample : PB123244BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB123244BL

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-BP091619.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	5.422	562	567	570	rBV	4602381	4293431	40.52%	7.195%
2	6.428	733	738	741	rBV	4485964	4325482	40.83%	7.248%
3	6.563	757	761	764	rBV	5237758	4694915	44.31%	7.867%
4	6.792	796	800	803	rBV	1869902	1409715	13.31%	2.362%
5	6.951	822	827	830	rBV	6580773	5608950	52.94%	9.399%
6	7.351	890	895	898	rBV	4155653	3982050	37.59%	6.673%
7	8.075	1014	1018	1021	rBV	1936810	1780973	16.81%	2.984%
8	9.151	1196	1201	1204	rBV	9399865	8999878	84.95%	15.081%
9	9.833	1313	1317	1320	rBV	2543315	2229577	21.04%	3.736%
10	10.628	1447	1452	1455	rBV	4438739	4277095	40.37%	7.167%
11	11.327	1567	1571	1574	rBV	2884164	2379644	22.46%	3.988%
12	12.933	1839	1844	1847	rBV	10532442	10594533	100.00%	17.754%
13	13.992	2020	2024	2027	rBV	2819608	2520301	23.79%	4.223%
14	15.474	2270	2276	2281	rBV	2059368	2579049	24.34%	4.322%

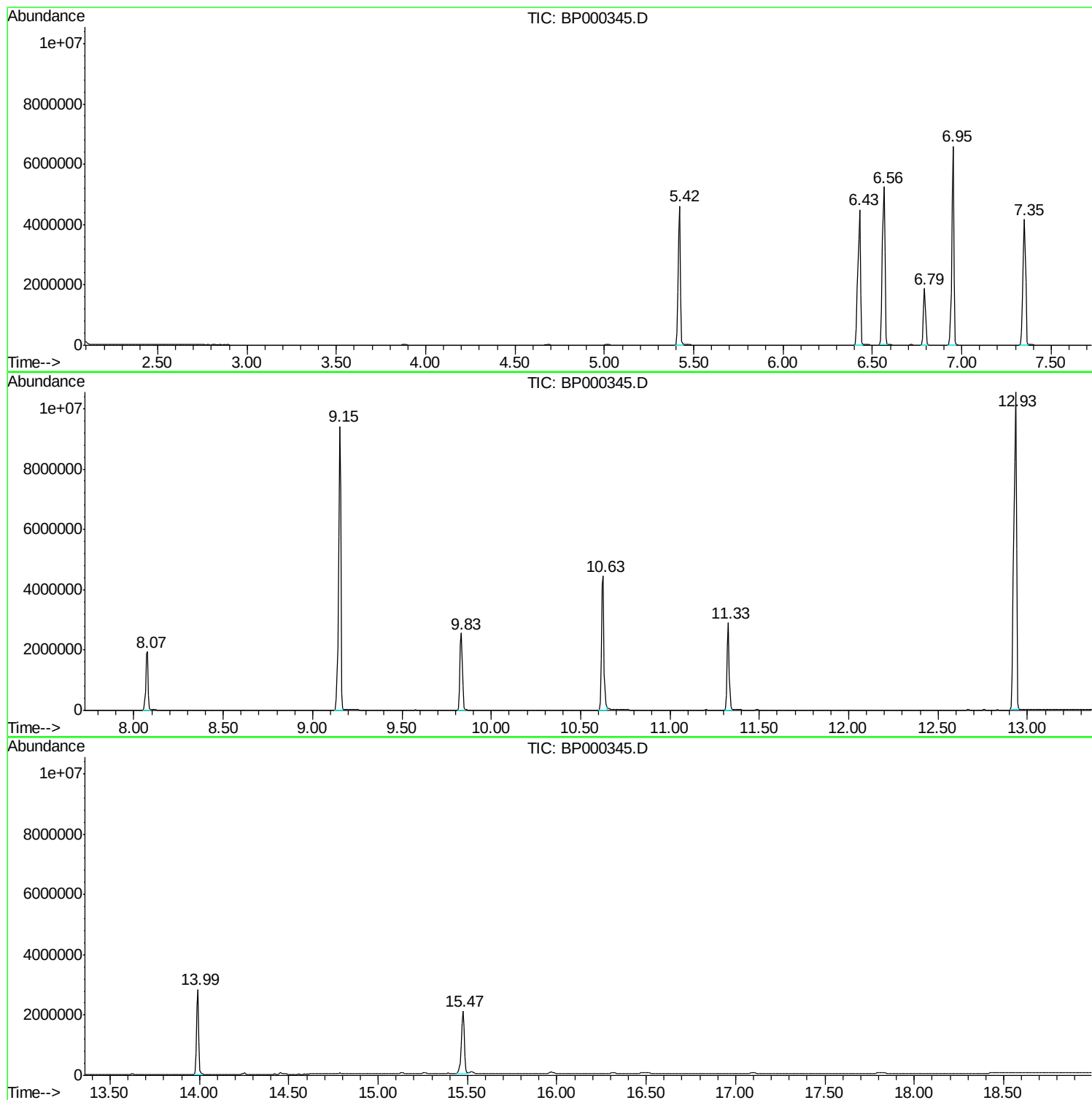
Sum of corrected areas: 59675593

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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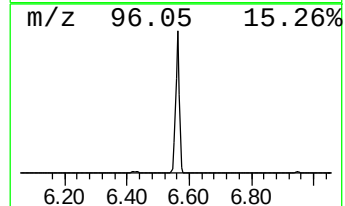
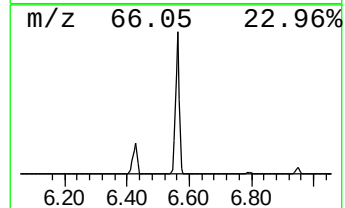
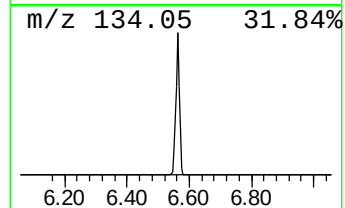
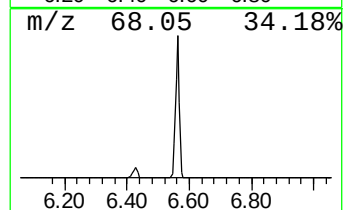
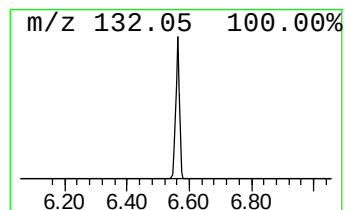
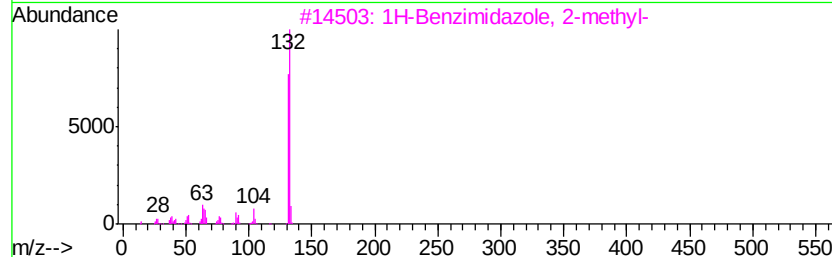
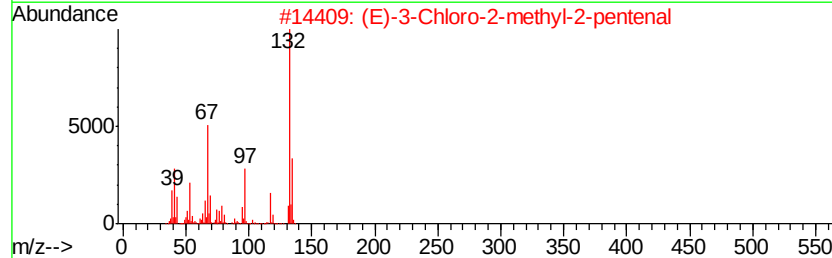
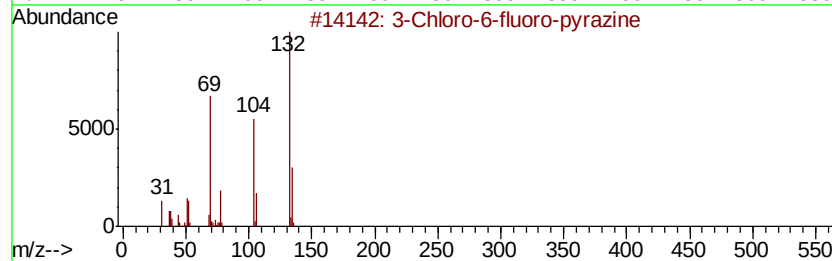
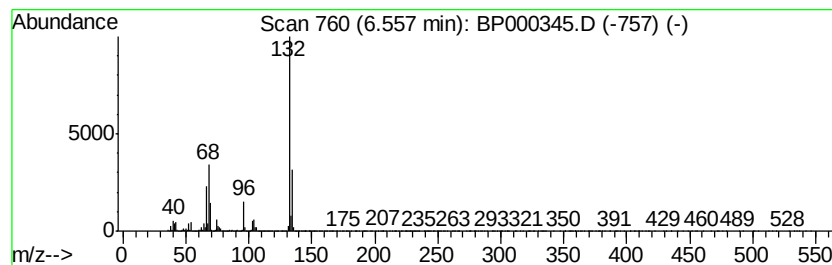
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	66.61 ng	4694920	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.56	6.56	66.6	ng	4694920	1	6.79	1409720	20.0