

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113569.D
 Acq On : 4 Apr 2019 3:49
 Operator : JU/SJ
 Sample : K2178-02 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-17

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.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	5.298	543	546	563	rBV	229996	310728	34.79%	4.107%
2	6.334	719	722	740	rBV	298120	397816	44.54%	5.258%
3	6.457	740	743	751	rVV	393379	367387	41.13%	4.856%
4	6.686	778	782	793	rBV	787733	712431	79.77%	9.416%
5	6.839	805	808	821	rBV	333892	306842	34.35%	4.055%
6	7.245	874	877	884	rBV	157777	190321	21.31%	2.515%
7	7.969	996	1000	1011	rBV	1019117	886949	99.30%	11.722%
8	9.045	1180	1183	1194	rBV	509194	431424	48.30%	5.702%
9	9.445	1247	1251	1259	rVB	21009	25499	2.85%	0.337%
10	9.722	1295	1298	1312	rVB2	1046020	893162	100.00%	11.804%
11	10.516	1430	1433	1443	rVB	151666	157365	17.62%	2.080%
12	11.204	1547	1550	1554	rBV	963806	864292	96.77%	11.423%
13	11.716	1632	1637	1639	rBV6	8565	12955	1.45%	0.171%
14	11.821	1652	1655	1658	rBV4	9431	13385	1.50%	0.177%
15	12.192	1714	1718	1722	rBV6	9640	22123	2.48%	0.292%
16	12.416	1753	1756	1759	rBV	62393	58378	6.54%	0.772%
17	12.645	1792	1795	1798	rBV	59311	51296	5.74%	0.678%
18	12.786	1815	1819	1823	rBV	501596	443306	49.63%	5.859%
19	13.845	1995	1999	2002	rBV	811702	777045	87.00%	10.270%
20	15.257	2235	2239	2246	rVB	488737	643665	72.07%	8.507%

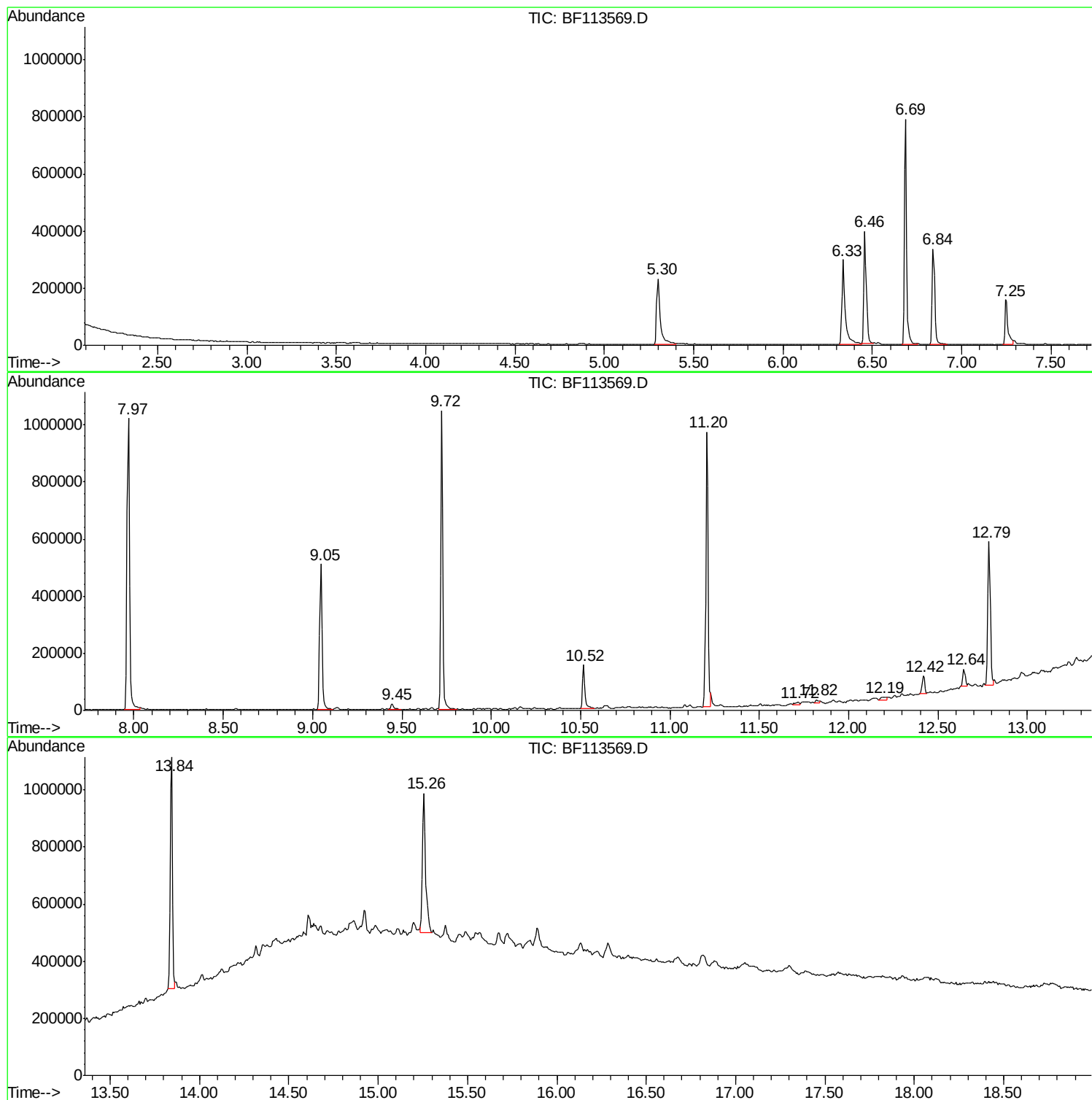
Sum of corrected areas: 7566369

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.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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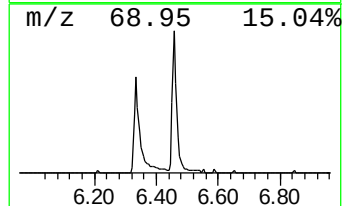
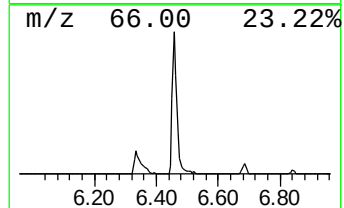
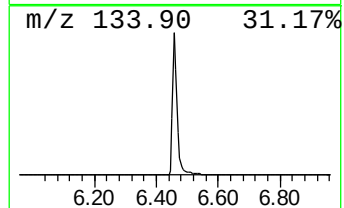
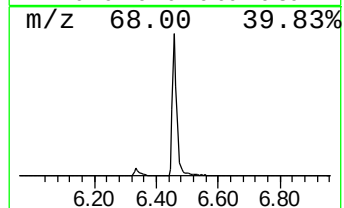
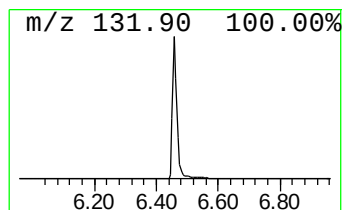
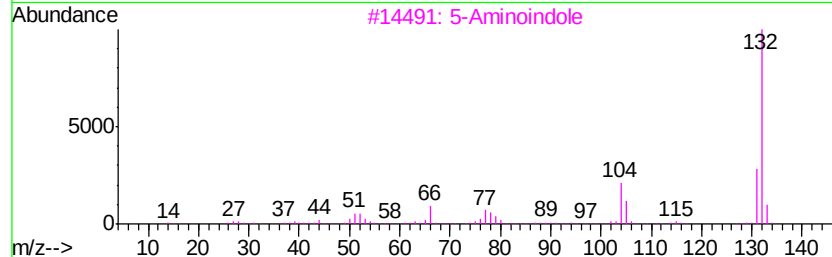
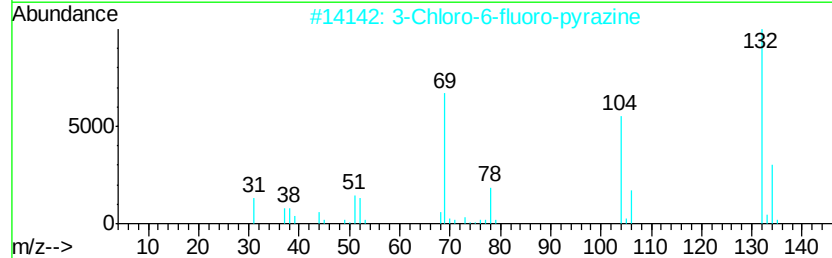
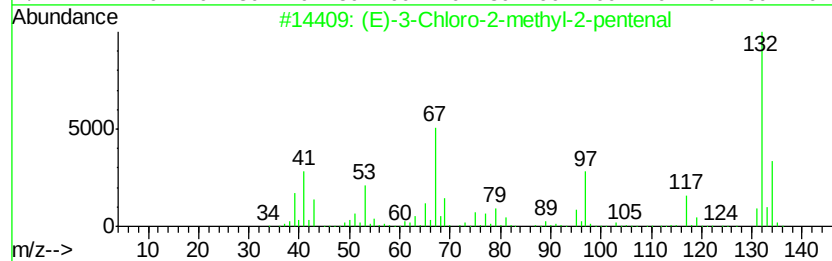
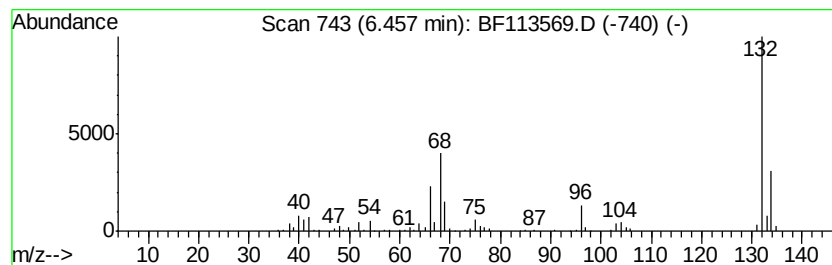
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	10.31 ng	367387	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Amphetaminil	250	C17H18N2	017590-01-1	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.46	6.46	10.3	ng	367387	1	6.69	712431	20.0