

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
 Data File : BF099968.D
 Acq On : 30 Oct 2017 19:03
 Operator : SJ/JU
 Sample : I6194-01 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-01-102717

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:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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	4.987	490	493	510	rBV	109289	144987	20.70%	1.936%
2	5.410	563	565	590	rBV	207366	332623	47.49%	4.441%
3	6.434	736	739	757	rBV	317325	479899	68.52%	6.408%
4	6.563	757	761	769	rVB	459076	467267	66.71%	6.239%
5	6.786	796	799	811	rBV	547892	475391	67.87%	6.348%
6	6.939	822	825	837	rBV	464116	402934	57.53%	5.380%
7	7.357	893	896	909	rBV	269392	284588	40.63%	3.800%
8	8.069	1014	1017	1030	rBV	814040	678683	96.90%	9.062%
9	8.633	1110	1113	1118	rBB2	9845	11359	1.62%	0.152%
10	9.139	1196	1199	1210	rBV	724723	596815	85.21%	7.969%
11	9.539	1264	1267	1275	rVB	91293	79529	11.35%	1.062%
12	9.686	1289	1292	1297	rBV	5080	8062	1.15%	0.108%
13	9.780	1304	1308	1312	rBV4	6047	12337	1.76%	0.165%
14	9.827	1312	1316	1320	rBV	790076	700412	100.00%	9.352%
15	10.539	1434	1437	1442	rBV	46973	42675	6.09%	0.570%
16	10.622	1448	1451	1460	rBV	112193	120702	17.23%	1.612%
17	10.727	1464	1469	1472	rBV2	8605	12533	1.79%	0.167%
18	11.321	1566	1570	1573	rBV	746703	645322	92.13%	8.617%
19	11.345	1573	1574	1578	rVV	60585	40630	5.80%	0.543%
20	11.398	1581	1583	1590	rVB	9775	13053	1.86%	0.174%
21	11.698	1631	1634	1637	rBV2	10804	8120	1.16%	0.108%
22	11.827	1653	1656	1658	rBV	6839	8083	1.15%	0.108%
23	11.857	1658	1661	1665	rVB	8302	10125	1.45%	0.135%
24	11.939	1672	1675	1677	rBV	16037	18486	2.64%	0.247%
25	12.380	1747	1750	1752	rBV2	11691	13993	2.00%	0.187%
26	12.486	1765	1768	1772	rVB4	11541	14339	2.05%	0.191%
27	12.539	1774	1777	1782	rVB	123928	111535	15.92%	1.489%
28	12.774	1810	1817	1820	rBV	105581	123778	17.67%	1.653%
29	12.904	1835	1839	1843	rVB	542633	450611	64.34%	6.017%
30	13.333	1909	1912	1917	rBV7	12927	20787	2.97%	0.278%
31	13.792	1986	1990	1994	rBV7	35154	50830	7.26%	0.679%
32	13.980	2017	2022	2025	rBV	523264	554254	79.13%	7.401%
33	14.004	2025	2026	2029	rVB	59618	40327	5.76%	0.538%
34	15.039	2199	2202	2204	rBV	52957	63729	9.10%	0.851%

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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	15.409	2262	2265	2267	rBV	30562	38765	5.53%	0.518%
36	15.468	2271	2275	2279	rVB	322590	411717	58.78%	5.497%

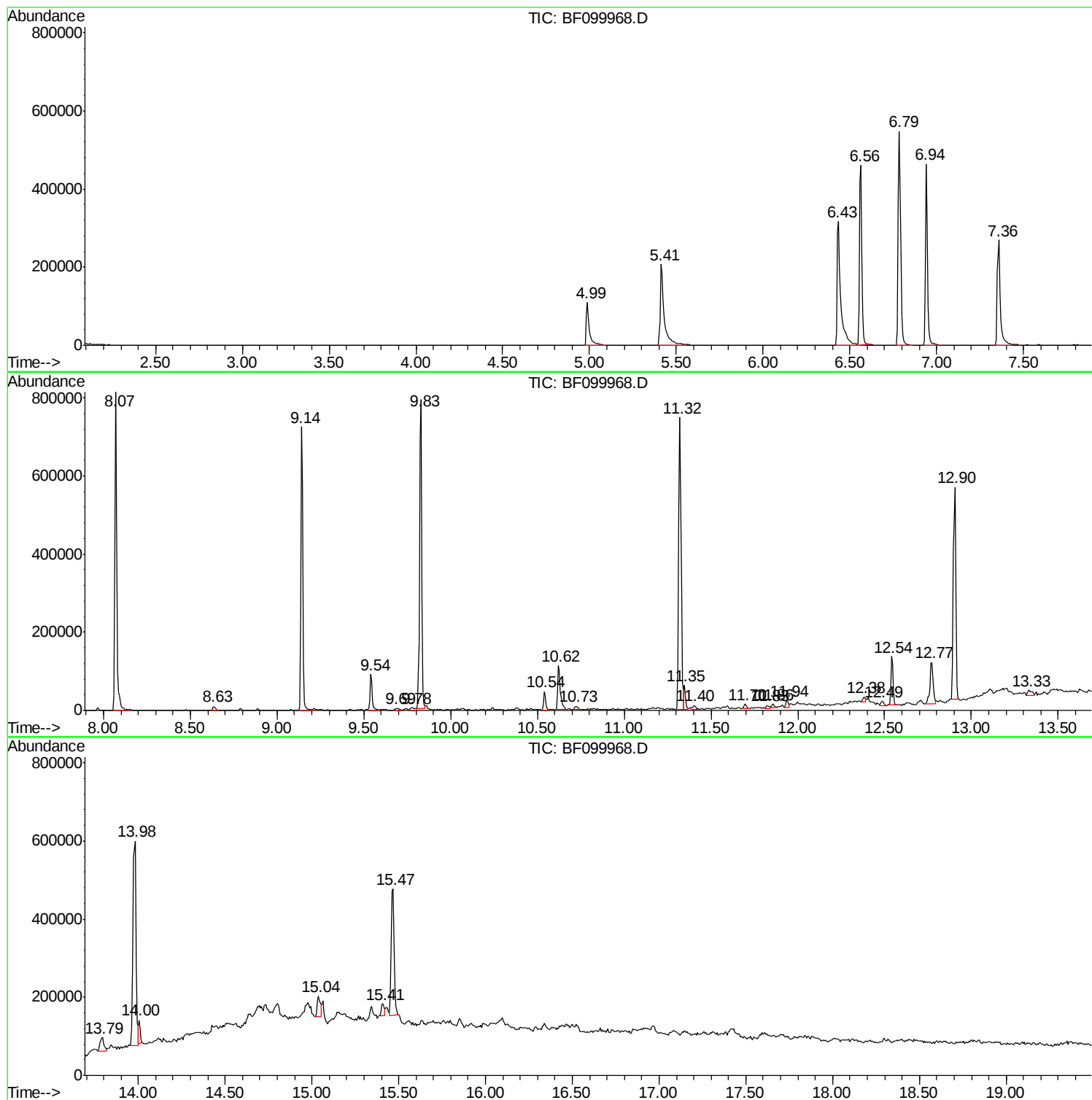
Sum of corrected areas: 7489280

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.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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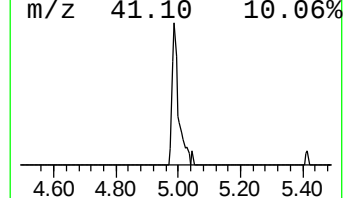
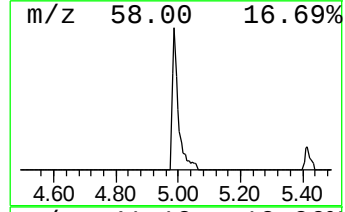
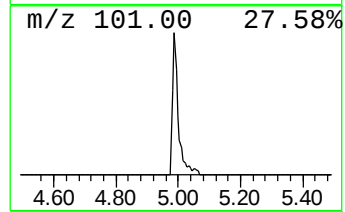
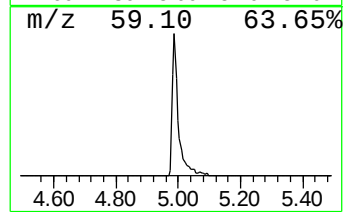
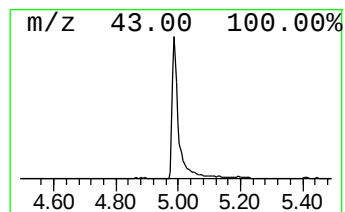
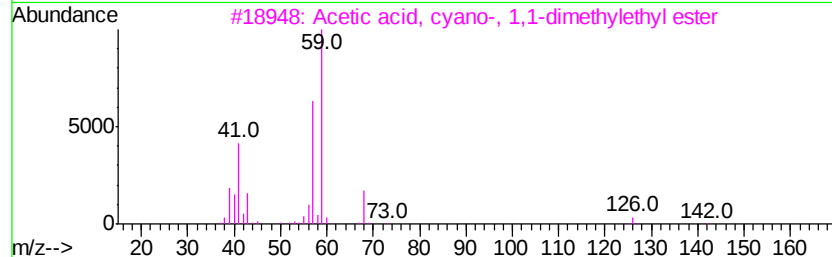
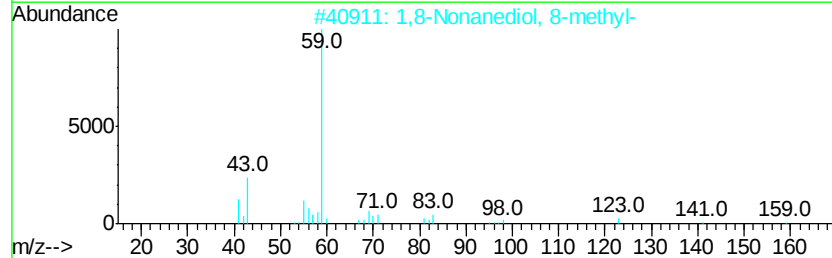
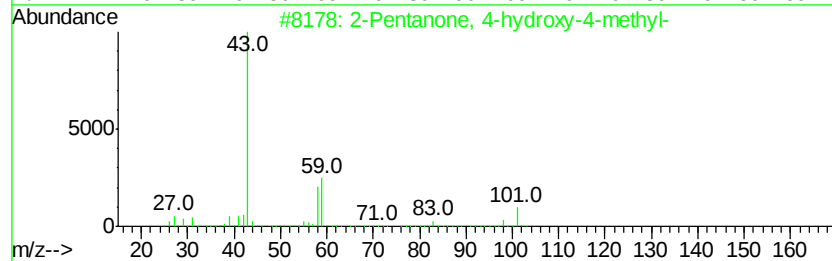
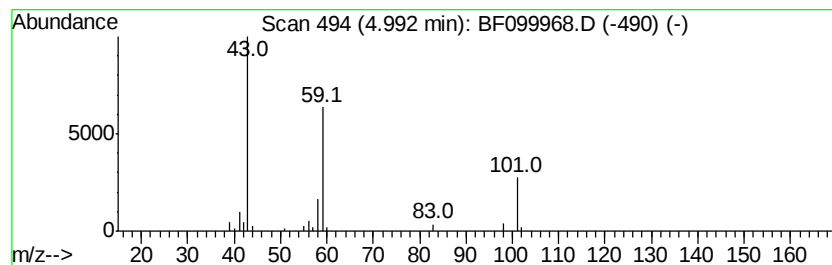
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	6.10 ng	144987	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	64
2			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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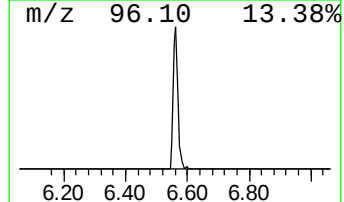
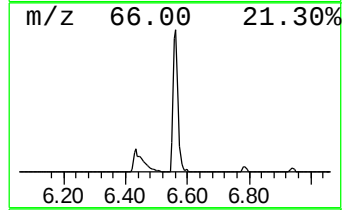
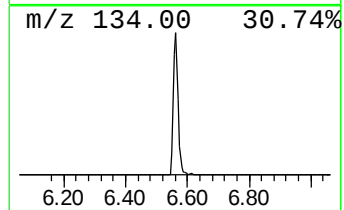
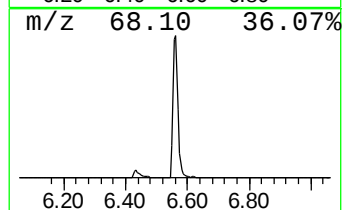
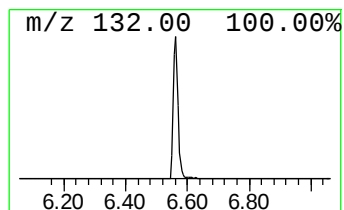
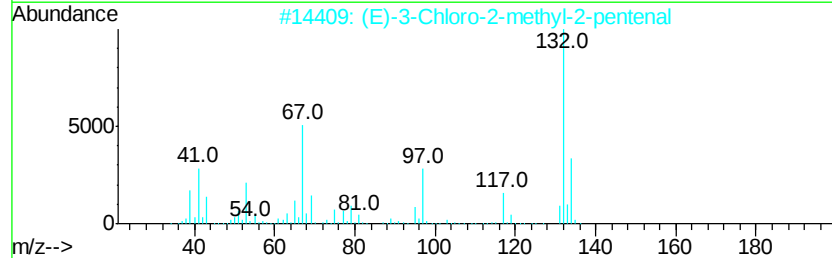
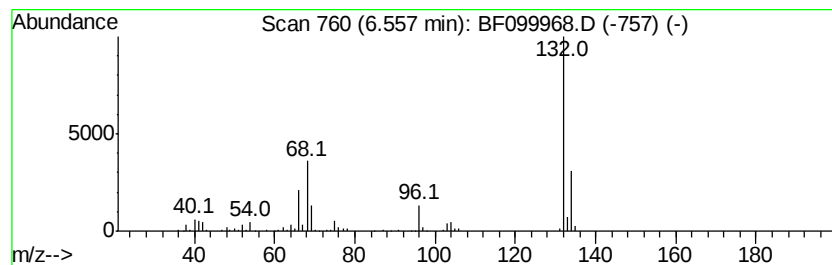
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	19.66 ng	467267	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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	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
2-Pentanone, 4-hy...	4.99	6.1	ng	144987	1	6.79	475391	20.0
unknown6.56	6.56	19.7	ng	467267	1	6.79	475391	20.0