

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070623\
 Data File : BF134118.D
 Acq On : 07 Jul 2023 03:02
 Operator : CG\JU
 Sample : 03376-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB-GW-062123

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_F\Methods\8270-BF062623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF134118.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.151	4	11	28	rVB	3696256	5095108	100.00%	24.910%
2	5.075	505	508	521	rVB	148715	149418	2.93%	0.731%
3	5.475	572	576	595	rBV	987694	909158	17.84%	4.445%
4	6.469	742	745	757	rBV	565540	526052	10.32%	2.572%
5	6.845	805	809	820	rBV	698082	633586	12.44%	3.098%
6	7.404	900	904	907	rBV	1808602	1640301	32.19%	8.020%
7	8.122	1023	1026	1042	rVB	1094855	894866	17.56%	4.375%
8	9.198	1205	1209	1213	rBV	3043966	2732310	53.63%	13.358%
9	9.880	1321	1325	1336	rBV	1058570	882077	17.31%	4.313%
10	10.674	1456	1460	1478	rBV	1959848	1774008	34.82%	8.673%
11	11.374	1575	1579	1587	rBV	1167118	1005447	19.73%	4.916%
12	12.974	1846	1851	1854	rBV	3340204	3008101	59.04%	14.707%
13	14.033	2027	2031	2041	rVB	691490	650786	12.77%	3.182%
14	15.539	2281	2287	2297	rBV	395158	552549	10.84%	2.701%

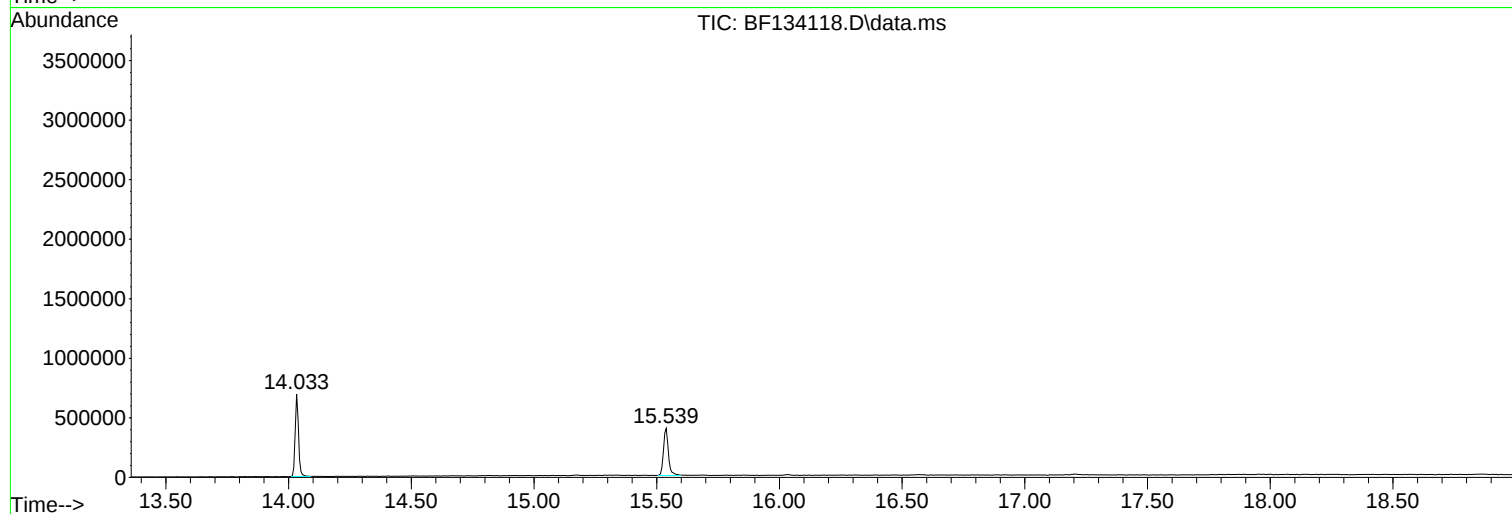
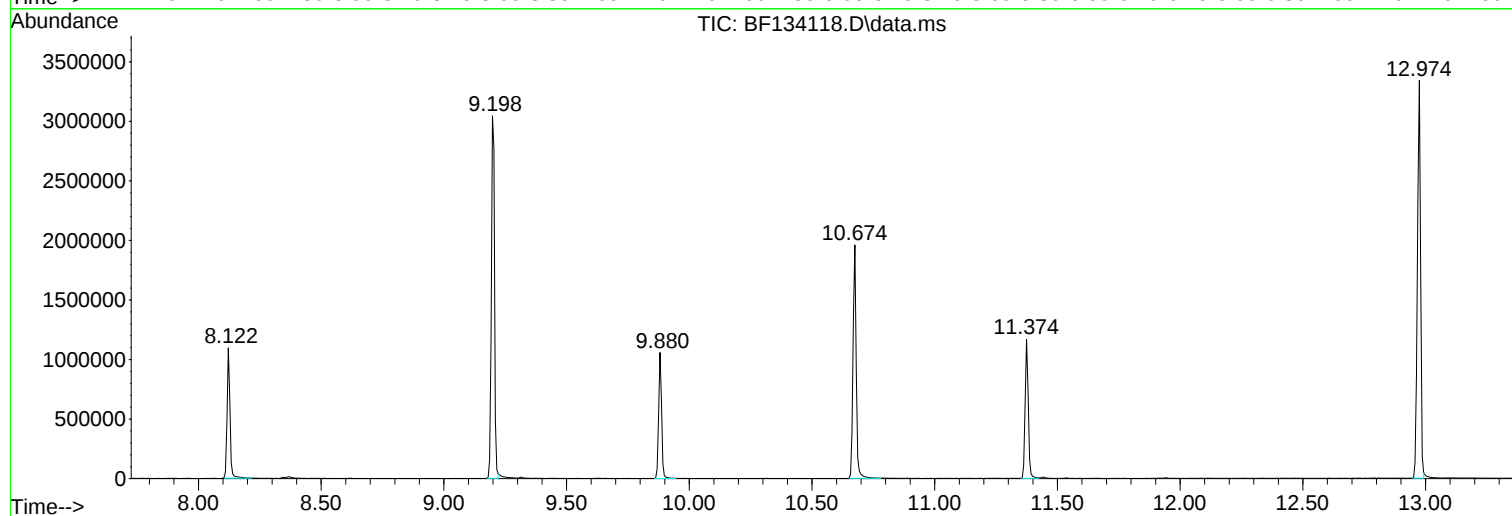
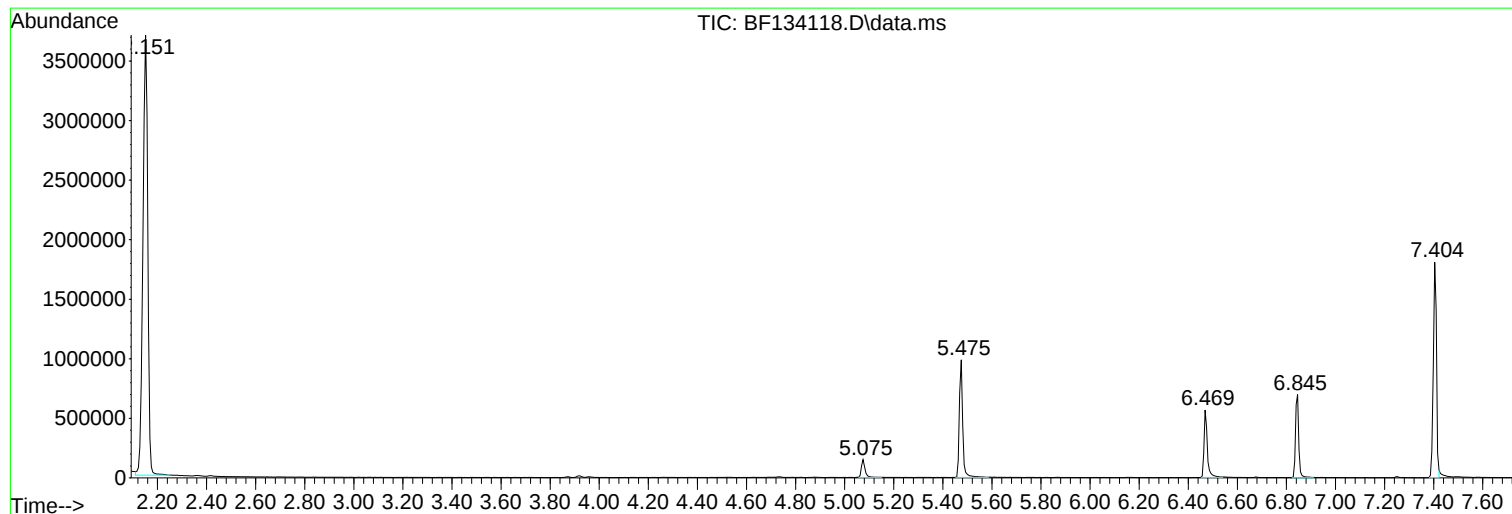
Sum of corrected areas: 20453767

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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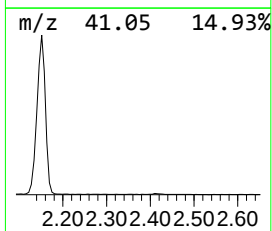
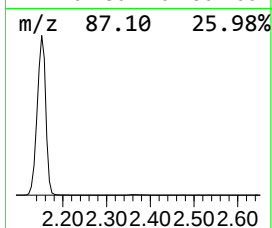
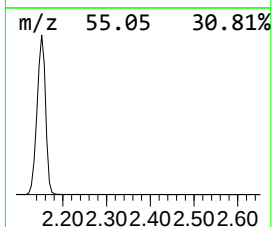
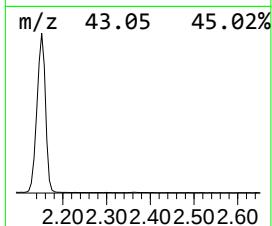
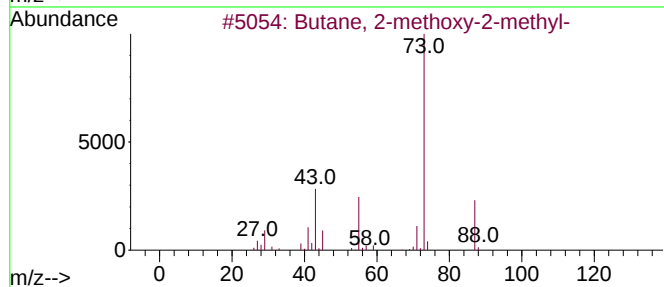
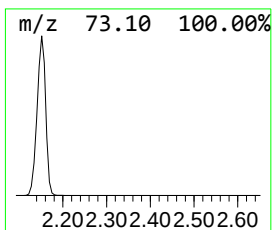
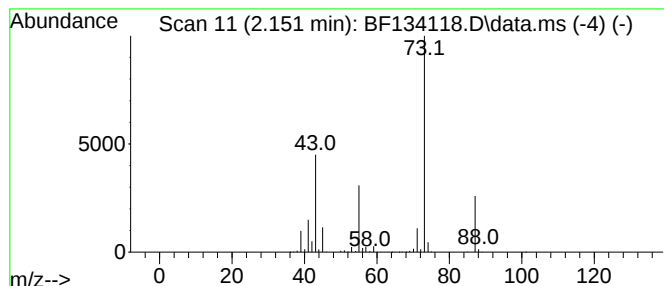
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.151	160.83 ng	5095110	1,4-Dichlorobenzene-d4	6.845

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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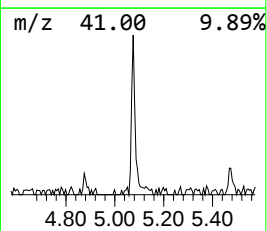
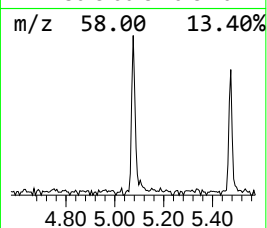
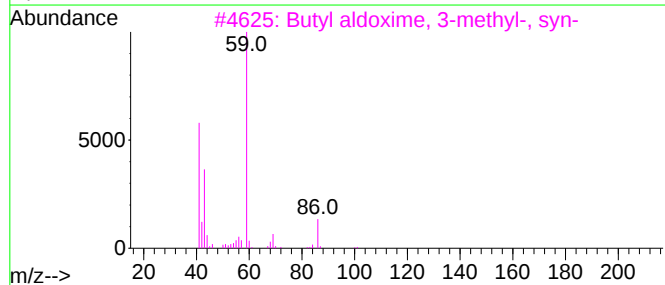
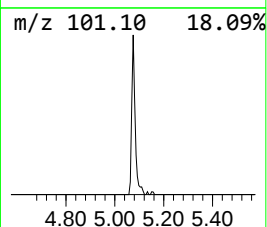
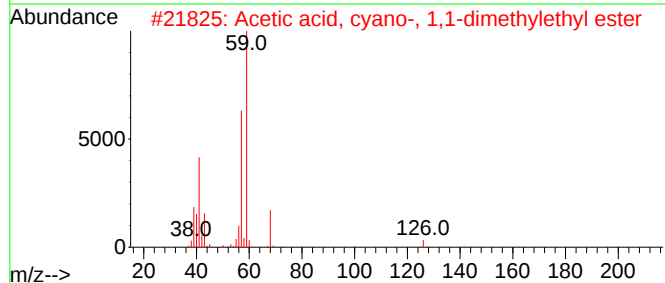
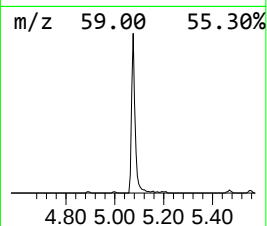
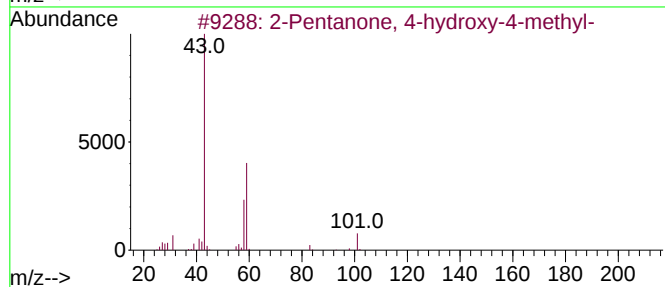
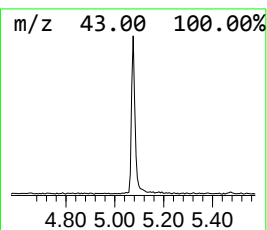
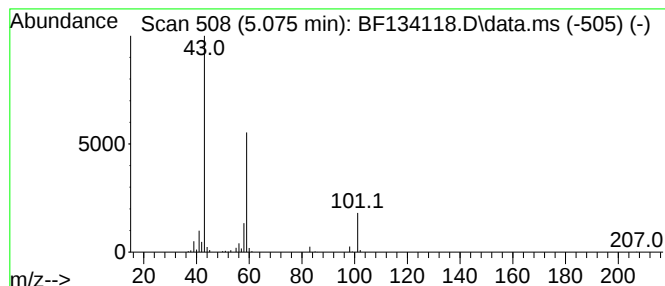
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.075 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	4.72 ng	149418	1,4-Dichlorobenzene-d4	6.845

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45	
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23	
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.151	160.8	ng	5095110	1	6.845	633586	20.0
unknown5.075	5.075	4.7	ng	149418	1	6.845	633586	20.0