

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081123\
 Data File : BF134731.D
 Acq On : 11 Aug 2023 14:13
 Operator : CG\JU
 Sample : PB154767BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB154767BL

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-BF080123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF134731.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.140	5	9	18	rVB	80470	119836	1.91%	0.314%
2	5.063	498	506	516	rBV	173876	265654	4.24%	0.697%
3	5.446	565	571	574	rBV	4529051	4895907	78.09%	12.847%
4	6.440	734	740	743	rBV	4184736	4881202	77.85%	12.809%
5	6.798	797	801	804	rBV	1445637	1130096	18.02%	2.966%
6	7.363	891	897	900	rBV	3065668	3280199	52.32%	8.608%
7	8.075	1013	1018	1022	rBV	1698261	1522689	24.29%	3.996%
8	9.157	1196	1202	1205	rBV	5557839	5611830	89.50%	14.726%
9	9.828	1311	1316	1320	rBV	1745887	1724484	27.50%	4.525%
10	10.622	1445	1451	1455	rBV	3587180	3759790	59.97%	9.866%
11	11.322	1566	1570	1573	rBV	2106853	1803861	28.77%	4.734%
12	12.922	1836	1842	1845	rBV	5927463	6269946	100.00%	16.453%
13	13.969	2016	2020	2035	rBV	1654630	1593934	25.42%	4.183%
14	15.439	2265	2270	2283	rBV	825279	1248588	19.91%	3.276%

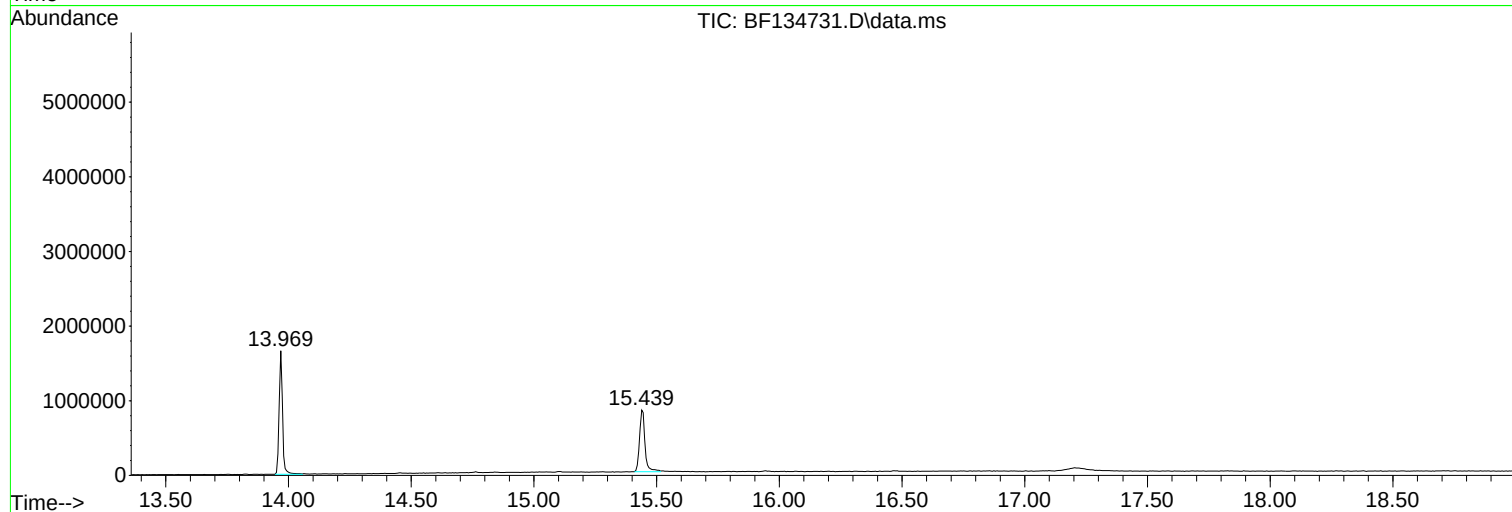
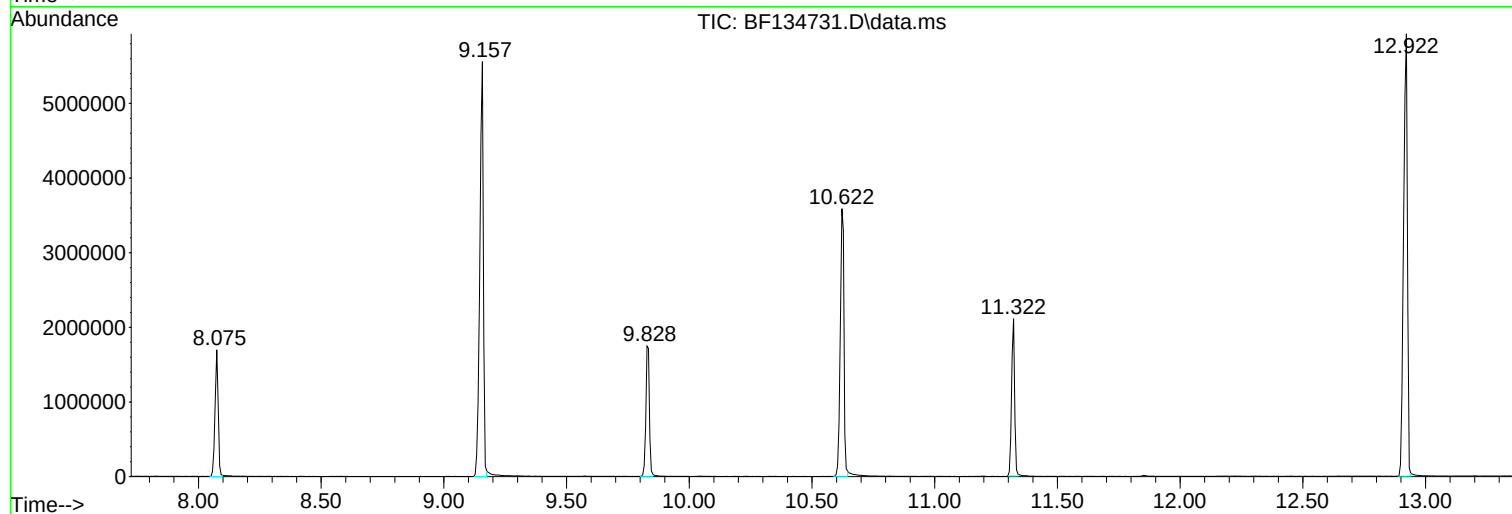
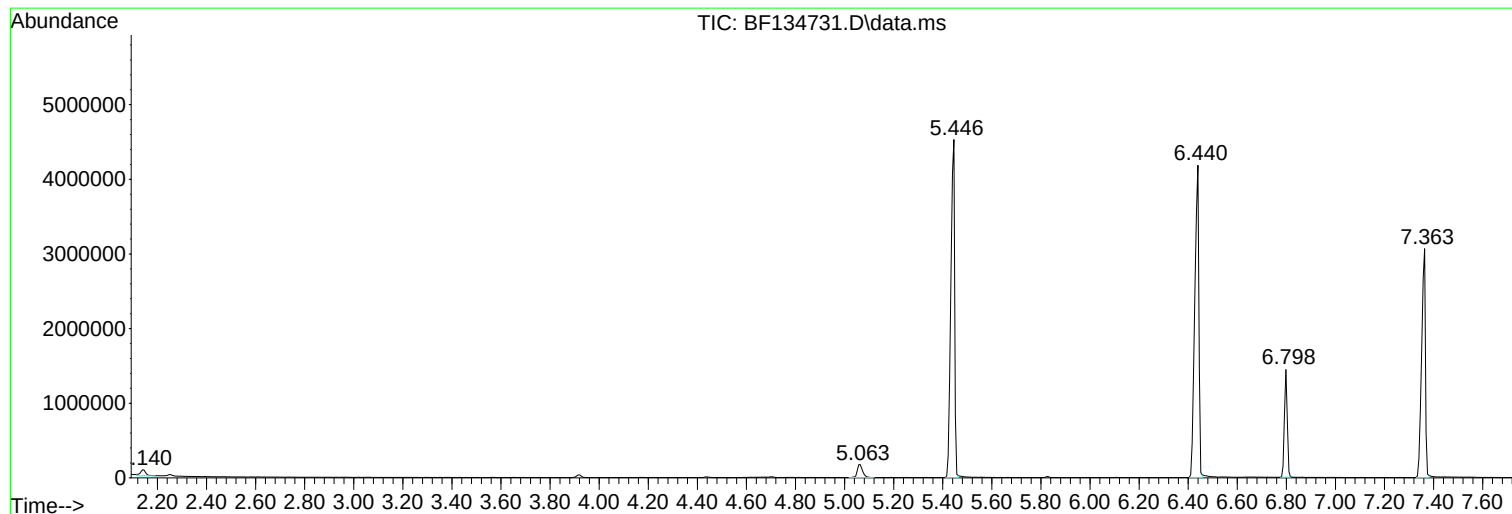
Sum of corrected areas: 38108016

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.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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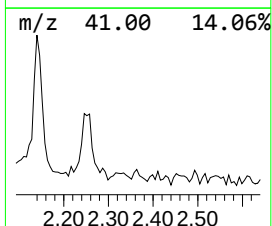
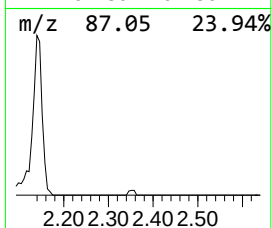
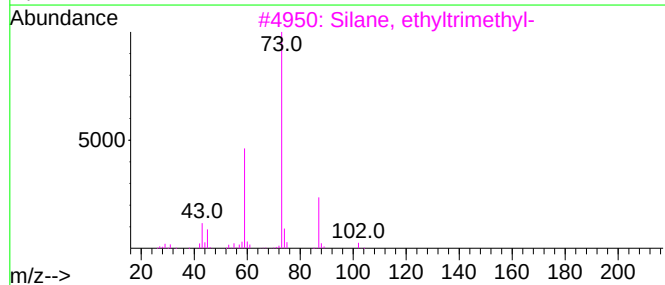
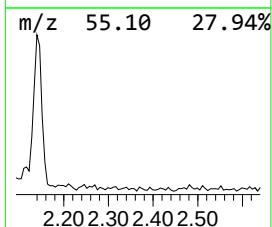
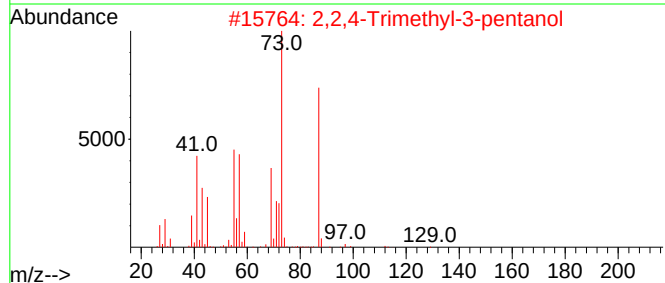
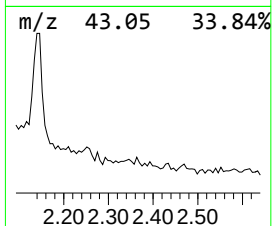
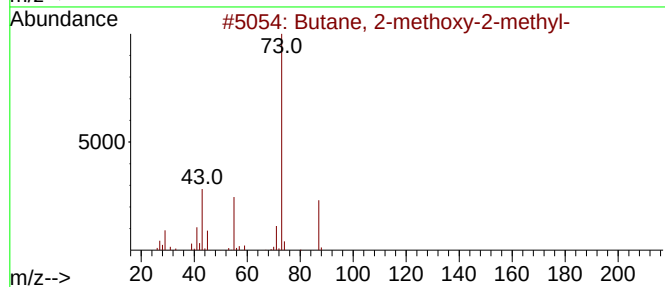
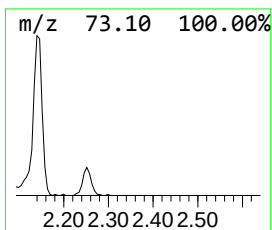
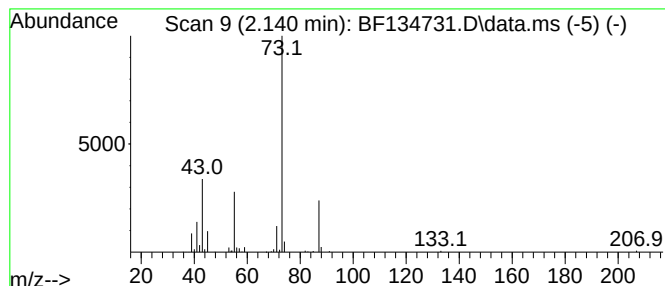
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.140	2.12 ng	119836	1,4-Dichlorobenzene-d4	6.798

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	40
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
4			2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	25
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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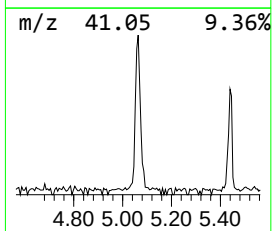
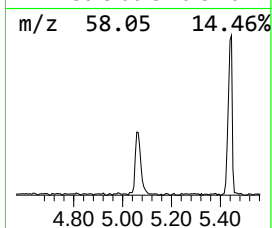
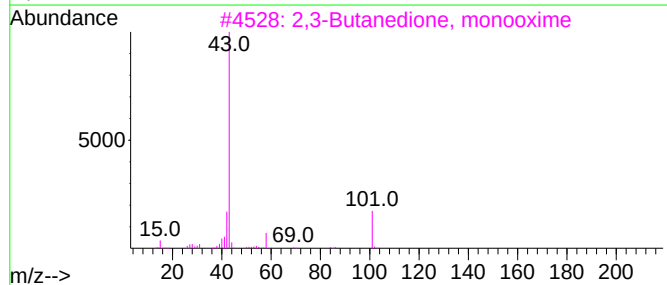
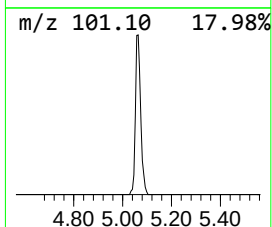
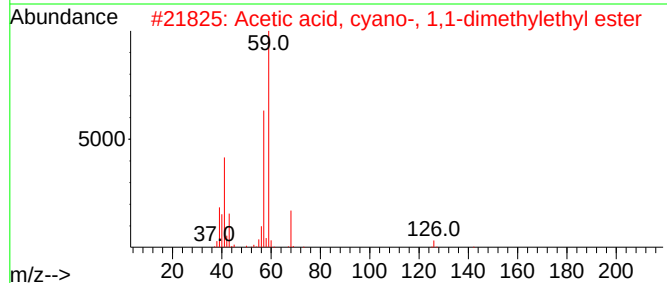
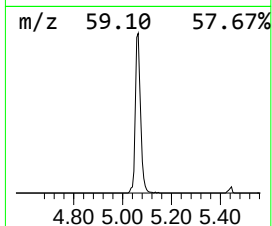
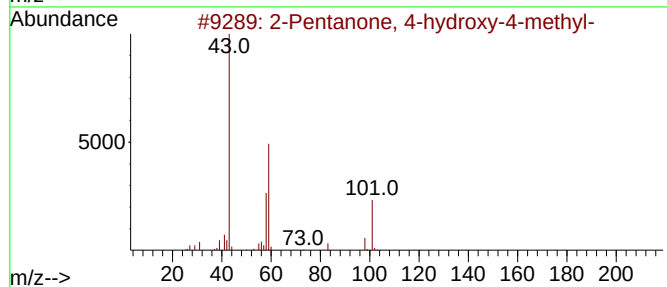
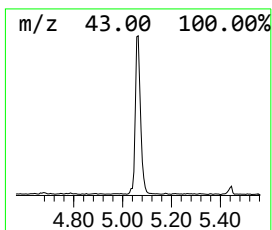
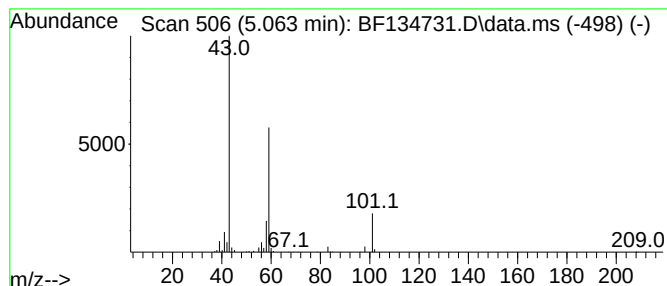
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	4.70 ng	265654	1,4-Dichlorobenzene-d4	6.798

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.140	2.1	ng	119836	1	6.798	1130100	20.0
2-Pentanone, 4-...	5.063	4.7	ng	265654	1	6.798	1130100	20.0