

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080625\
 Data File : BF143324.D
 Acq On : 06 Aug 2025 12:05
 Operator : RC/JU
 Sample : PB169119BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB169119BL

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-BF071725.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143324.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.357	17	28	36	rVB	35098	62096	1.51%	0.264%
2	5.216	506	514	531	rBV	248867	416486	10.13%	1.768%
3	5.604	574	580	598	rBV	2425115	3426793	83.37%	14.547%
4	6.592	741	748	753	rBV	2302137	3399584	82.71%	14.432%
5	6.963	806	811	816	rBV	534426	666956	16.23%	2.831%
6	7.516	898	905	911	rBV	1621263	2368187	57.62%	10.053%
7	8.239	1022	1028	1046	rBV	672484	884684	21.52%	3.756%
8	9.316	1204	1211	1243	rBV	2964415	4110329	100.00%	17.449%
9	9.998	1321	1327	1346	rBV	725890	946338	23.02%	4.017%
10	10.786	1455	1461	1480	rBV	1492065	2059669	50.11%	8.744%
11	11.486	1574	1580	1599	rBV	662371	873999	21.26%	3.710%
12	13.068	1843	1849	1869	rBV	2511113	3378198	82.19%	14.341%
13	14.127	2023	2029	2040	rBV	347926	501077	12.19%	2.127%
14	15.639	2280	2286	2303	rVB	230212	461998	11.24%	1.961%

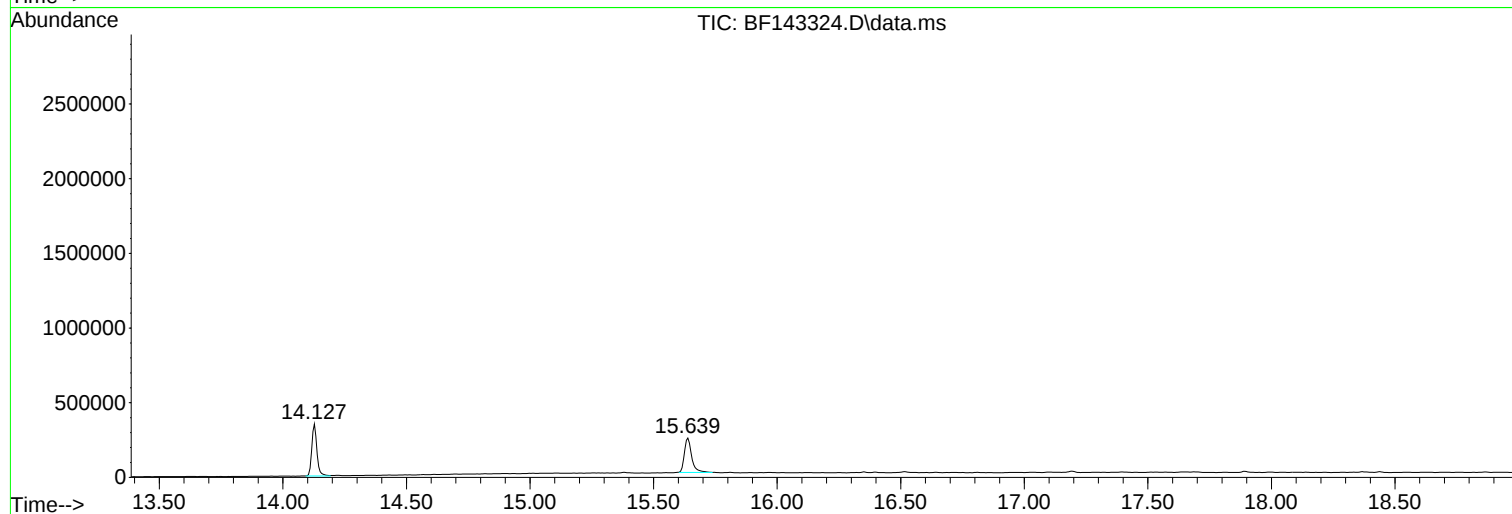
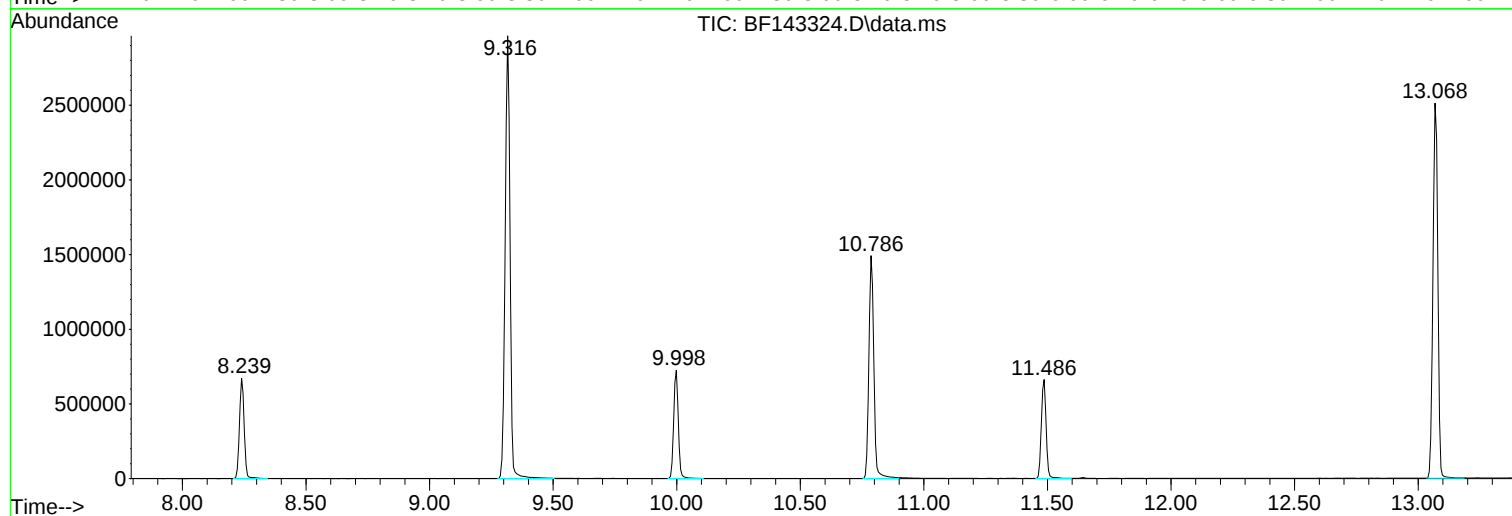
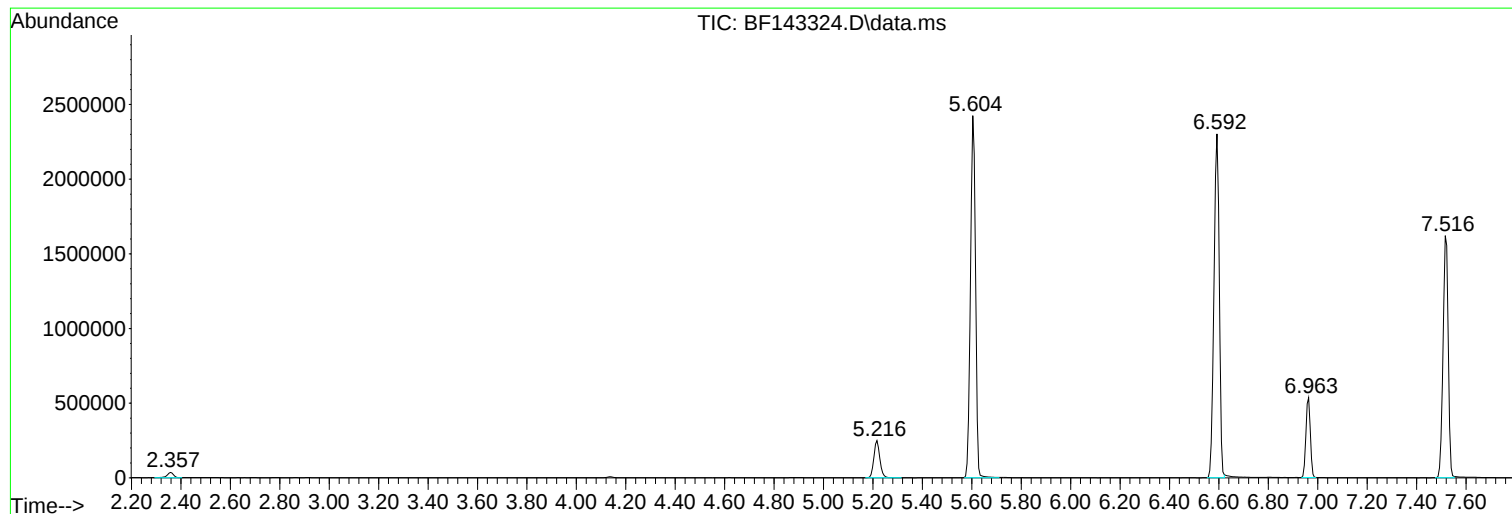
Sum of corrected areas: 23556394

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.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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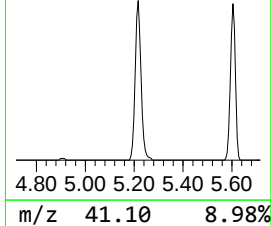
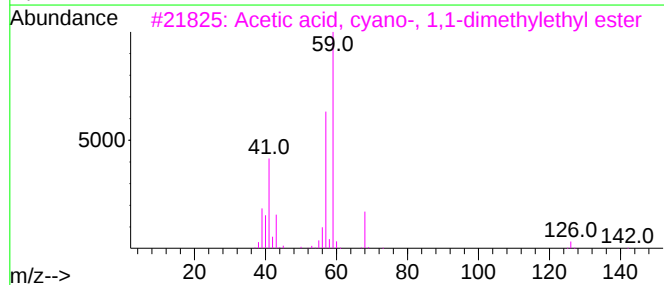
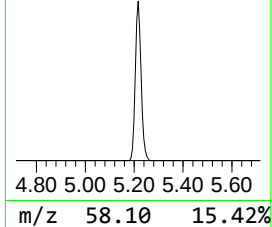
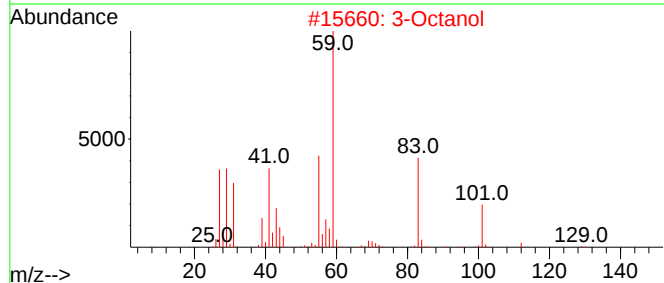
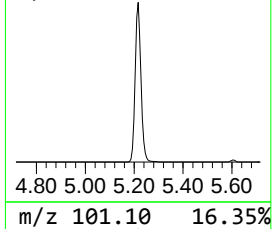
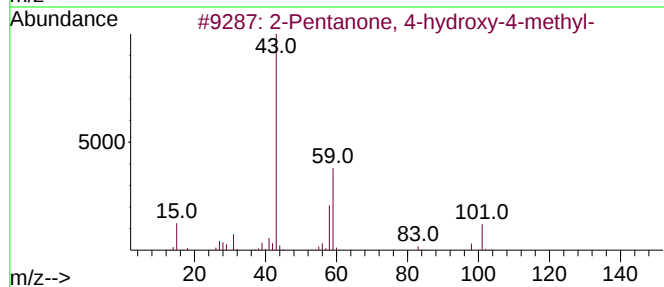
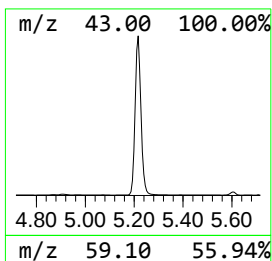
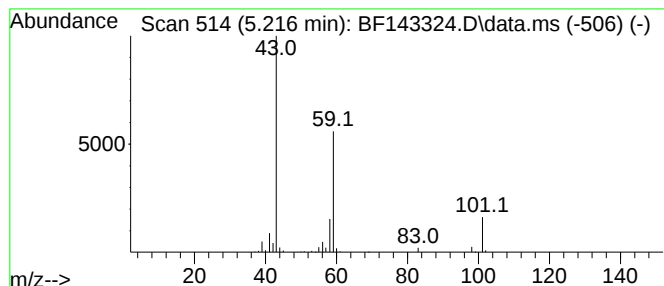
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.216	12.49 ng	416486	1,4-Dichlorobenzene-d4	6.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Octanol	130	C8H18O	000589-98-0	33		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.216	12.5	ng	416486	1	6.963	666956	20.0