

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091522\
 Data File : BP011720.D
 Acq On : 15 Sep 2022 23:57
 Operator : CG/JU
 Sample : PB147206BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147206BL

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_P\Methods\8270E-BP091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP011720.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.076	11	15	19	rBV	287248	337923	1.72%	0.353%
2	5.028	342	347	360	rBV	492263	745495	3.79%	0.778%
3	5.493	419	426	442	rBV	7258110	10896994	55.45%	11.377%
4	7.093	688	698	717	rBV	6807251	10926202	55.60%	11.407%
5	7.934	834	841	850	rBV	1472866	2332148	11.87%	2.435%
6	9.099	1031	1039	1053	rBV	3984264	6628431	33.73%	6.920%
7	10.746	1310	1319	1335	rBV	1903452	3248106	16.53%	3.391%
8	13.199	1728	1736	1751	rBV	9690302	14704966	74.82%	15.352%
9	14.575	1963	1970	1983	rBV	2615452	3896750	19.83%	4.068%
10	16.063	2216	2223	2245	rBV	6578451	9088290	46.24%	9.488%
11	17.328	2431	2438	2449	rBV	3193720	4363359	22.20%	4.555%
12	19.881	2866	2872	2877	rBV	16221915	19652732	100.00%	20.518%
13	21.404	3126	3131	3141	rBV2	3539316	4631372	23.57%	4.835%
14	23.857	3542	3548	3559	rVB	2257105	4331143	22.04%	4.522%

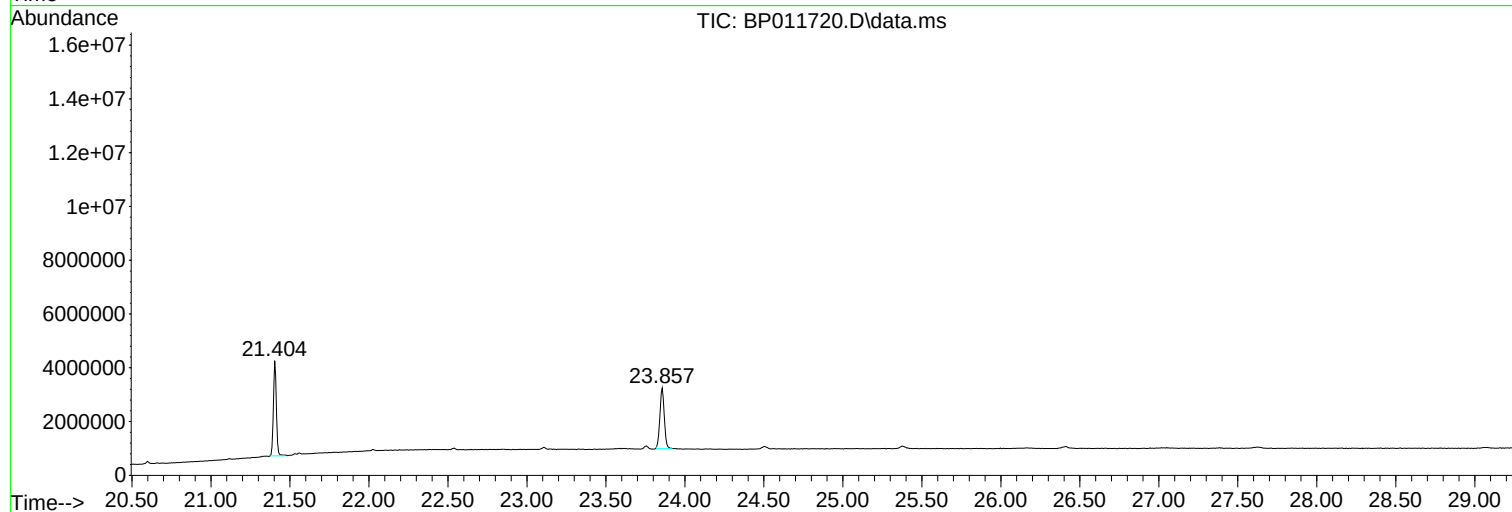
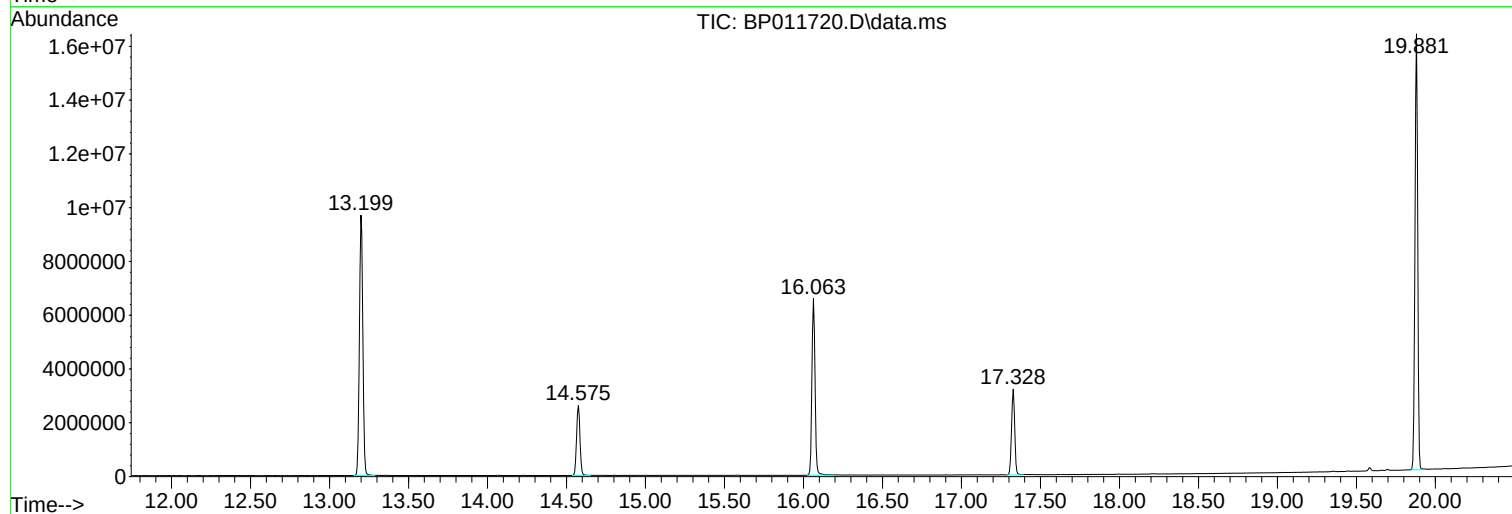
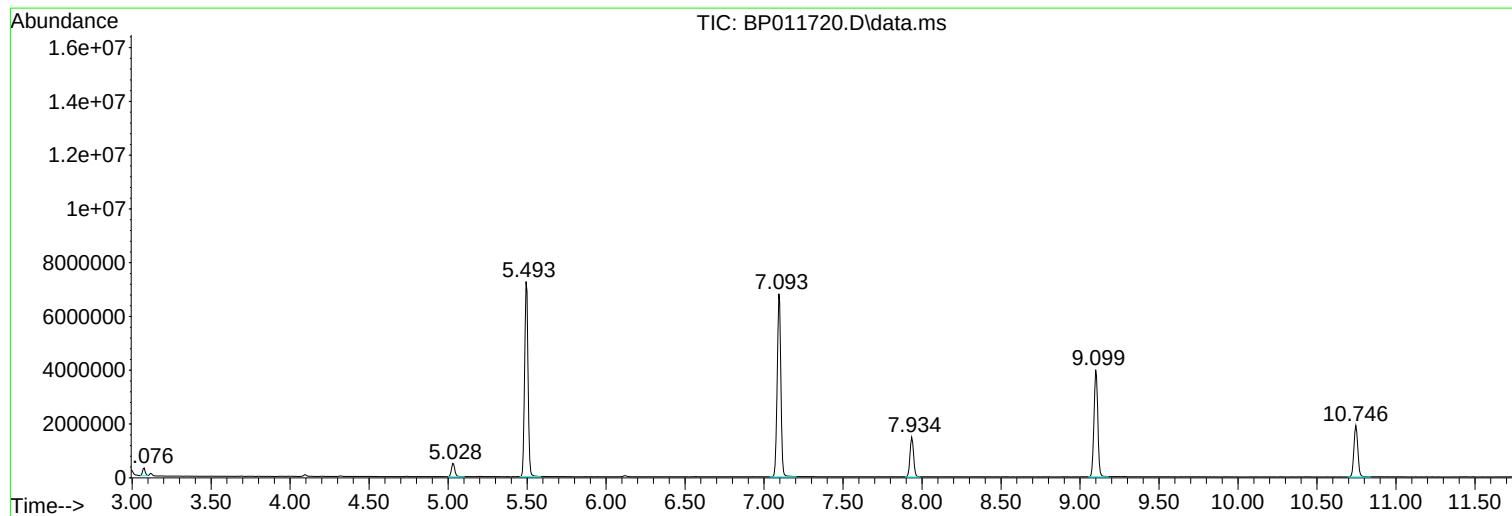
Sum of corrected areas: 95783911

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.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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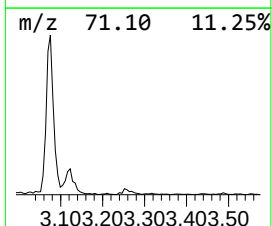
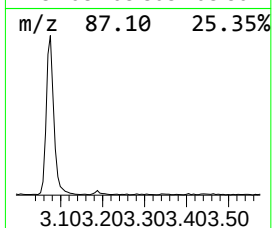
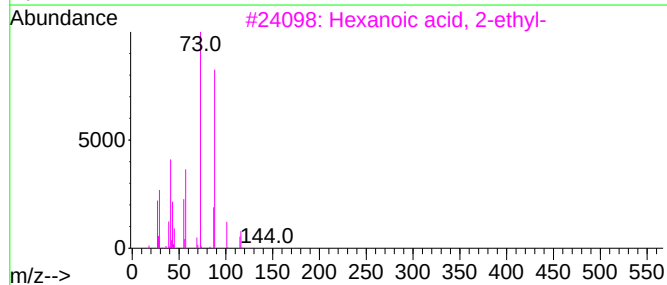
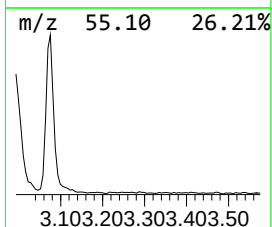
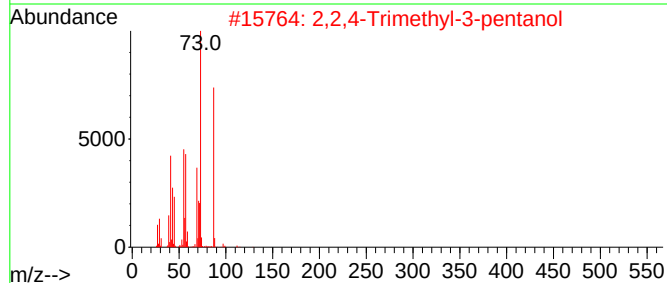
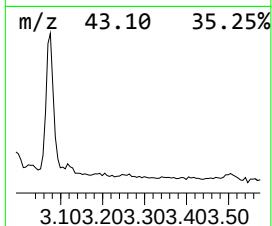
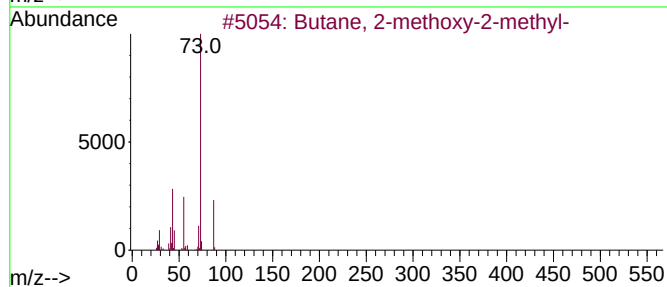
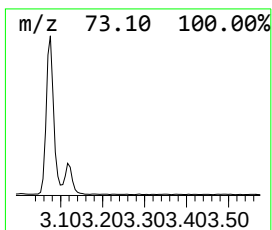
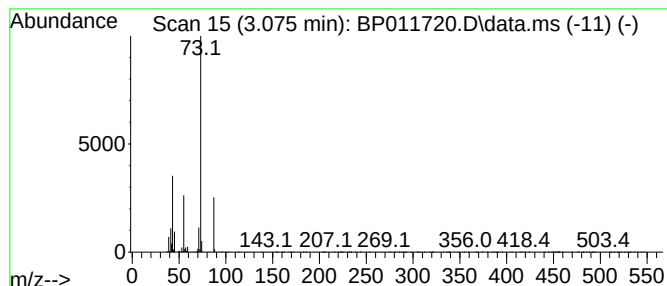
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	2.90 ng	337923	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	36
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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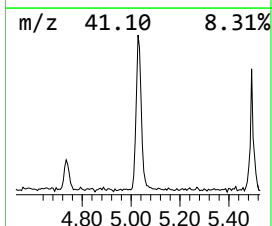
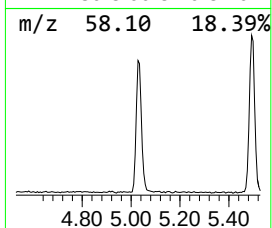
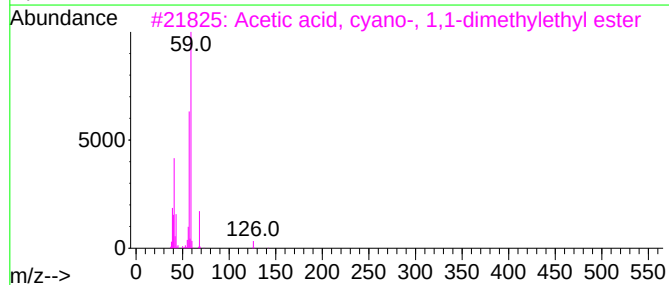
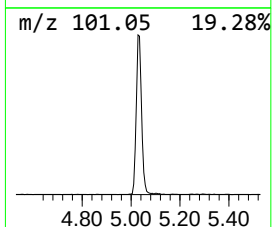
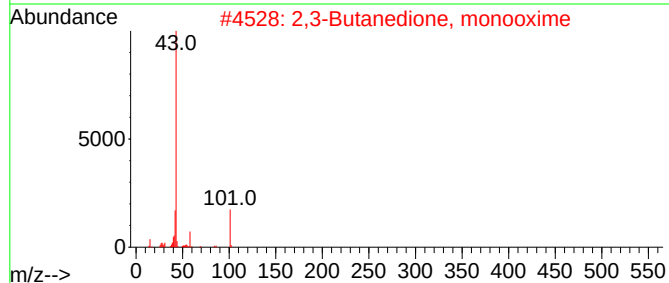
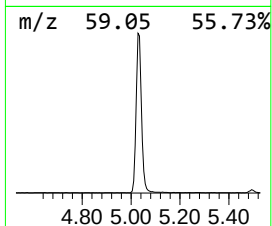
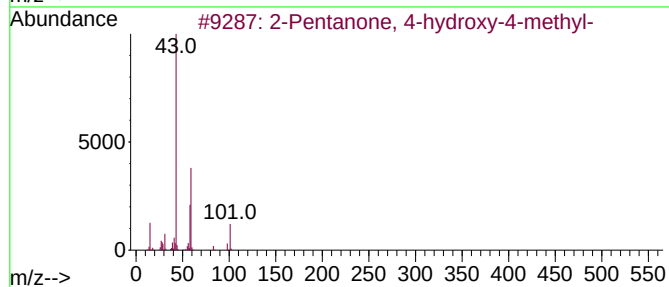
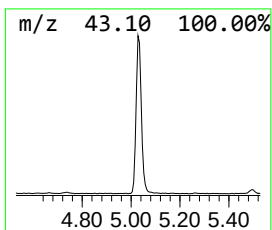
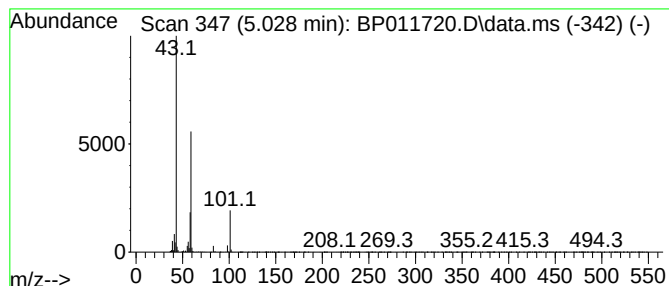
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.028	6.39 ng	745495	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Acetone	58	C3H6O	000067-64-1	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.075	2.9	ng	337923	1	7.934	2332150	20.0
2-Pentanone, 4-...	5.028	6.4	ng	745495	1	7.934	2332150	20.0