

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110724\
Data File : BE101525.D
Acq On : 7 Nov 2024 11:26
Operator : RC/JU
Sample : PB164711BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
PB164711BL

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:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BE101525.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.684	12	16	28	rVB	36116	58498	1.19%	0.280%
2	4.781	368	373	392	rBV	215965	347414	7.09%	1.663%
3	5.322	458	465	489	rBV	1078402	1835753	37.44%	8.786%
4	6.949	734	742	773	rBV	901440	1846964	37.67%	8.840%
5	7.566	839	847	855	rBV	239750	392183	8.00%	1.877%
6	8.776	1044	1053	1075	rBV	678636	1174756	23.96%	5.623%
7	10.333	1308	1318	1334	rBV	306663	551904	11.26%	2.642%
8	12.795	1729	1737	1756	rBV	1859207	2806764	57.24%	13.434%
9	14.176	1963	1972	1981	rBV	473436	732072	14.93%	3.504%
10	15.680	2221	2228	2249	rBV	1606531	2463109	50.23%	11.789%
11	16.914	2431	2438	2450	rBV	610138	893847	18.23%	4.278%
12	19.505	2873	2879	2884	rBV	3643572	4903303	100.00%	23.468%
13	21.074	3141	3146	3153	rBV	983144	1257201	25.64%	6.017%
14	23.359	3529	3535	3544	rBV2	816174	1629478	33.23%	7.799%

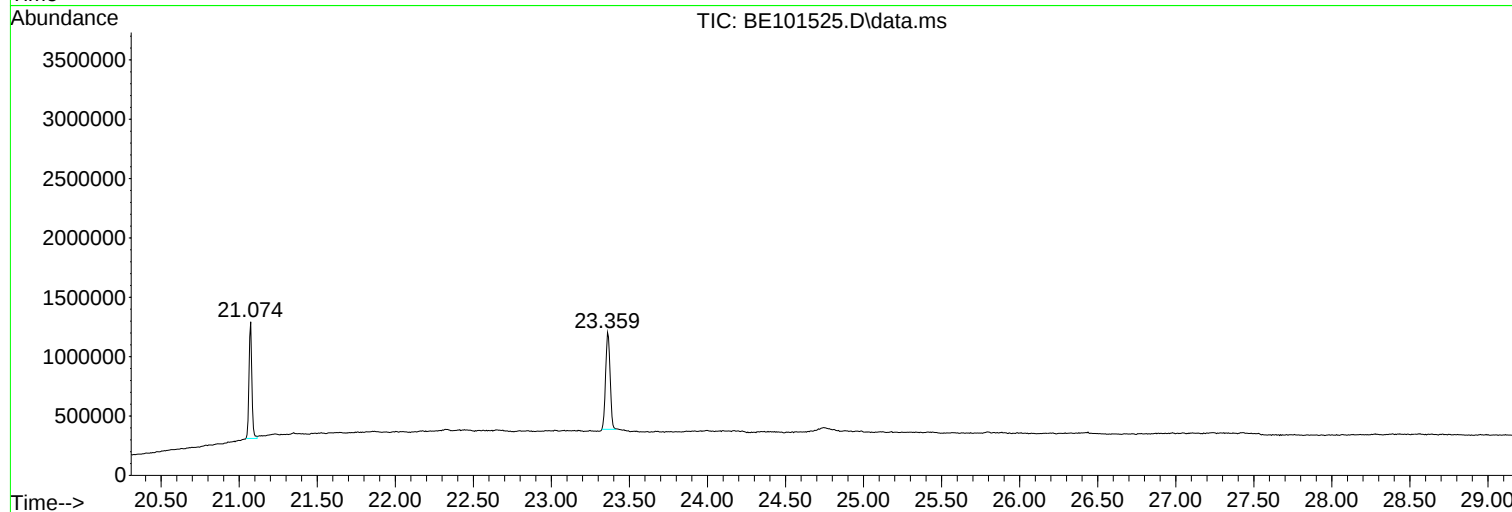
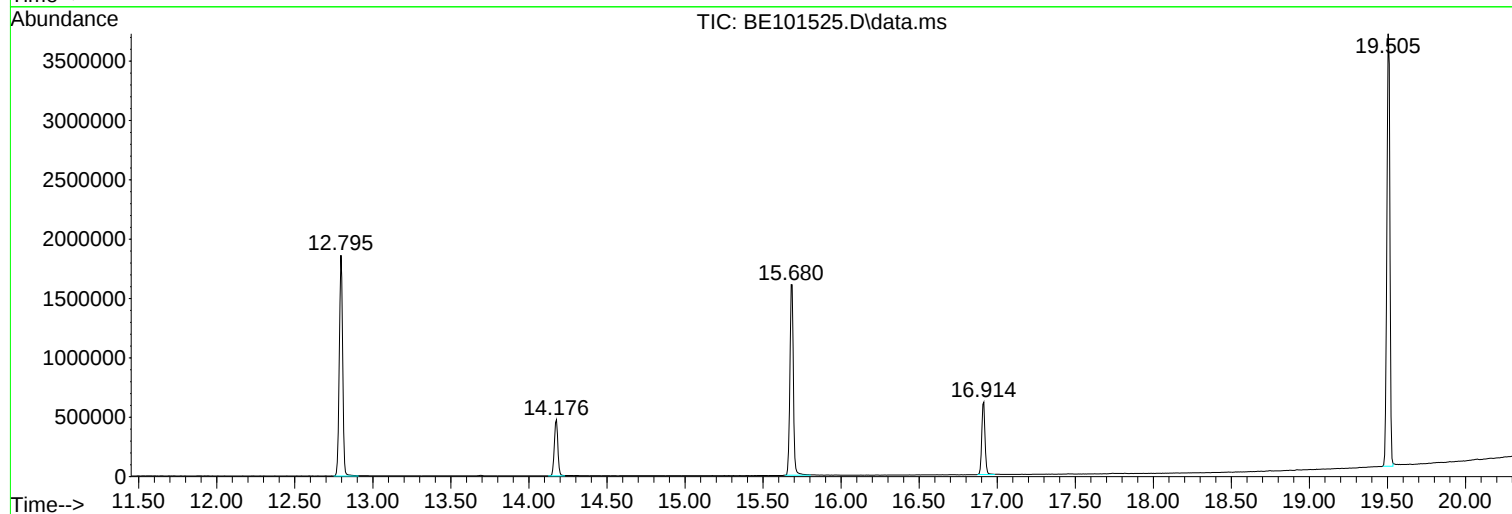
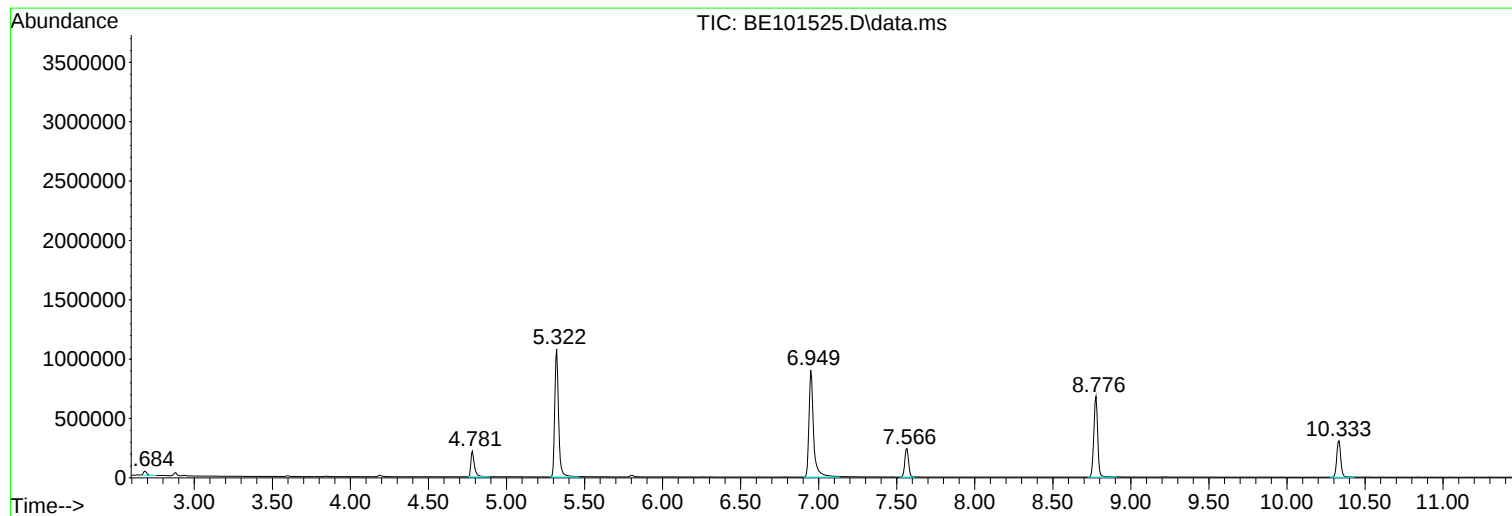
Sum of corrected areas: 20893246

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.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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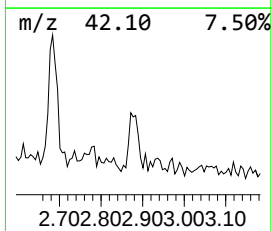
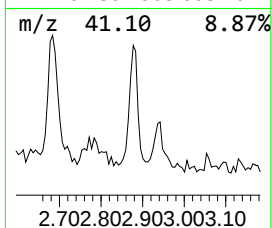
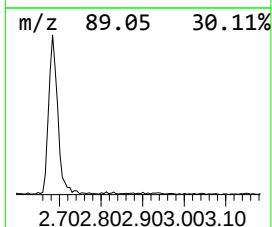
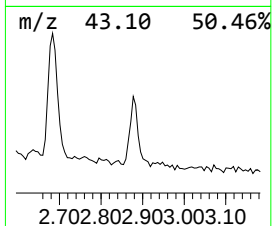
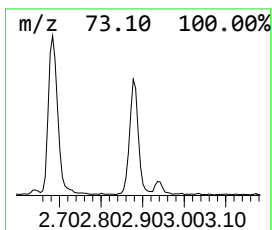
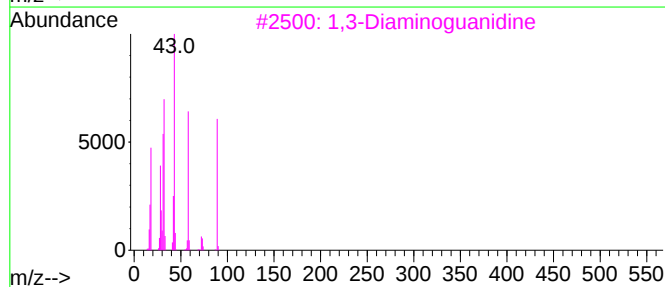
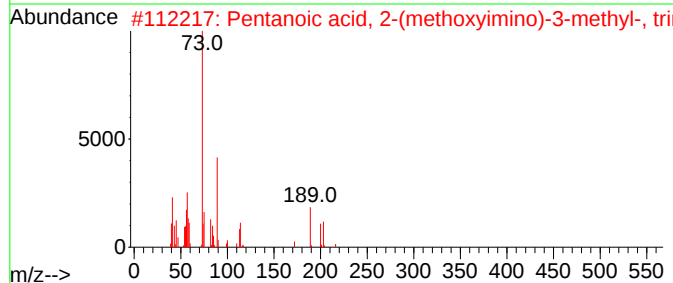
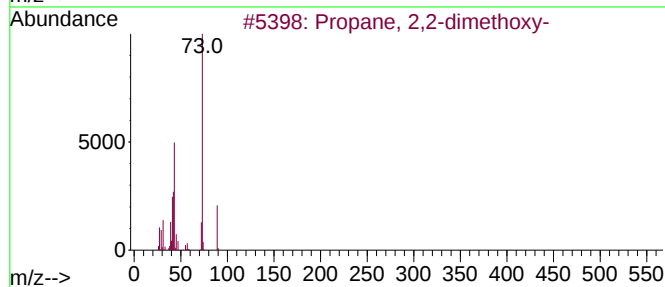
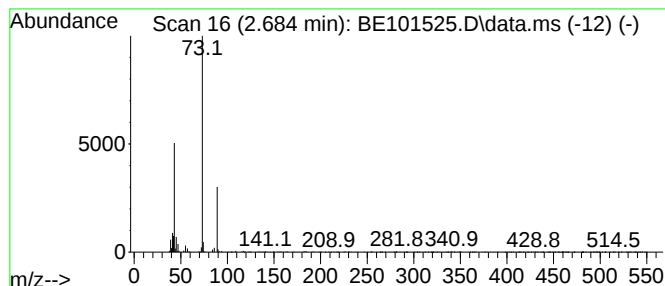
Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.684 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.684	2.98 ng	58498	1,4-Dichlorobenzene-d4	7.566

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	42
2			Pentanoic acid, 2-(methoxyimino)-...	231	C10H21NO3Si	055494-02-5	9
3			1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1,3-Dioxolane, 2-(6-octynyl)-	182	C11H18O2	056741-65-2	9



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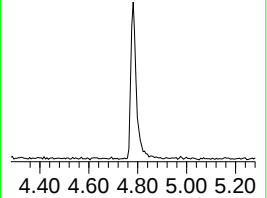
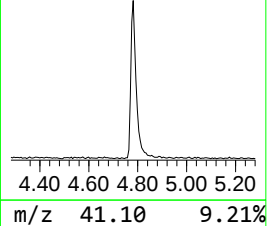
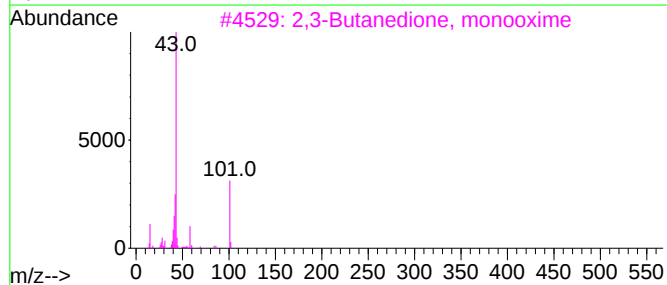
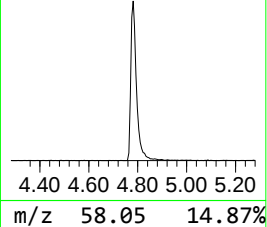
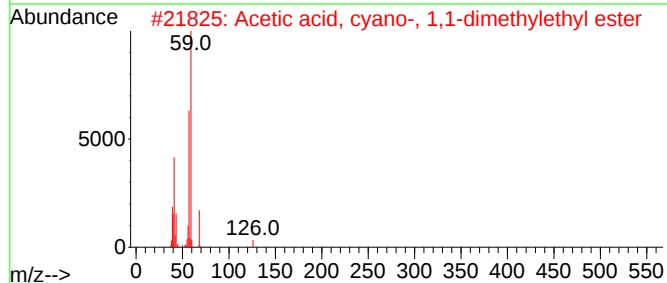
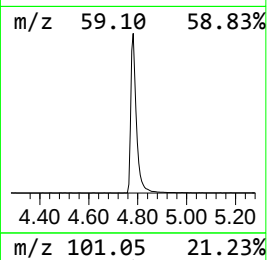
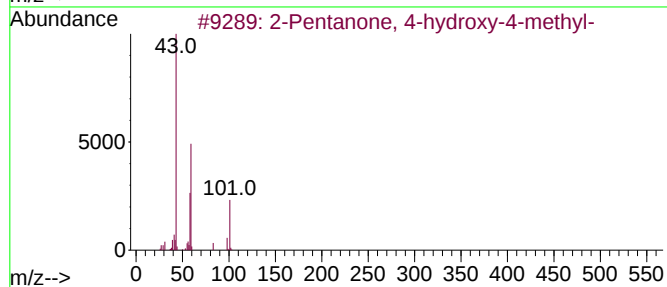
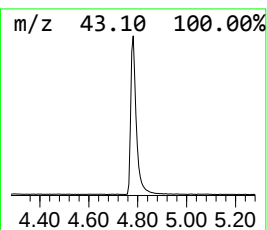
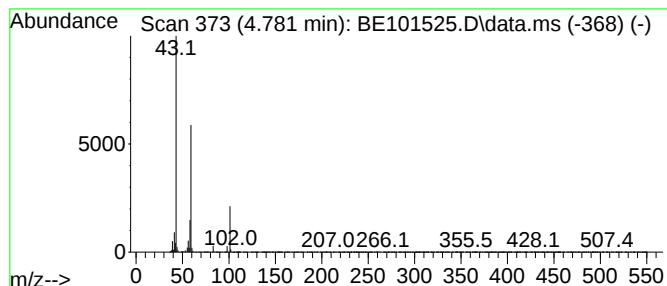
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.781	17.72 ng	347414	1,4-Dichlorobenzene-d4	7.566	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
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
unknown2.684	2.684	3.0	ng	58498	1	7.566	392183	20.0
2-Pentanone, 4-...	4.781	17.7	ng	347414	1	7.566	392183	20.0