

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110824\
 Data File : BP022937.D
 Acq On : 08 Nov 2024 16:51
 Operator : RC/JU
 Sample : PB164369BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB164369BL

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-BP110724.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP022937.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.781	299	305	316	rBV	61702	87704	1.41%	0.401%
2	5.234	375	382	409	rBV	1279598	1838041	29.54%	8.396%
3	6.740	631	638	641	rBV	52669	77774	1.25%	0.355%
4	6.793	641	647	672	rVB	1113009	1856661	29.84%	8.481%
5	7.593	776	783	794	rBV	228692	373285	6.00%	1.705%
6	8.740	967	978	993	rBV	681757	1207471	19.41%	5.516%
7	10.351	1242	1252	1267	rBV	256803	462006	7.43%	2.110%
8	12.816	1664	1671	1691	rBV	2102061	3001824	48.25%	13.712%
9	14.216	1903	1909	1923	rBV	435969	661167	10.63%	3.020%
10	15.728	2159	2166	2185	rBV	1760685	2693940	43.30%	12.306%
11	17.033	2379	2388	2402	rBV	511281	877897	14.11%	4.010%
12	19.757	2846	2851	2861	rBV	4967346	6221943	100.00%	28.422%
13	21.468	3135	3142	3151	rBV	736279	1236882	19.88%	5.650%
14	24.704	3684	3692	3704	rBV	482471	1294734	20.81%	5.914%

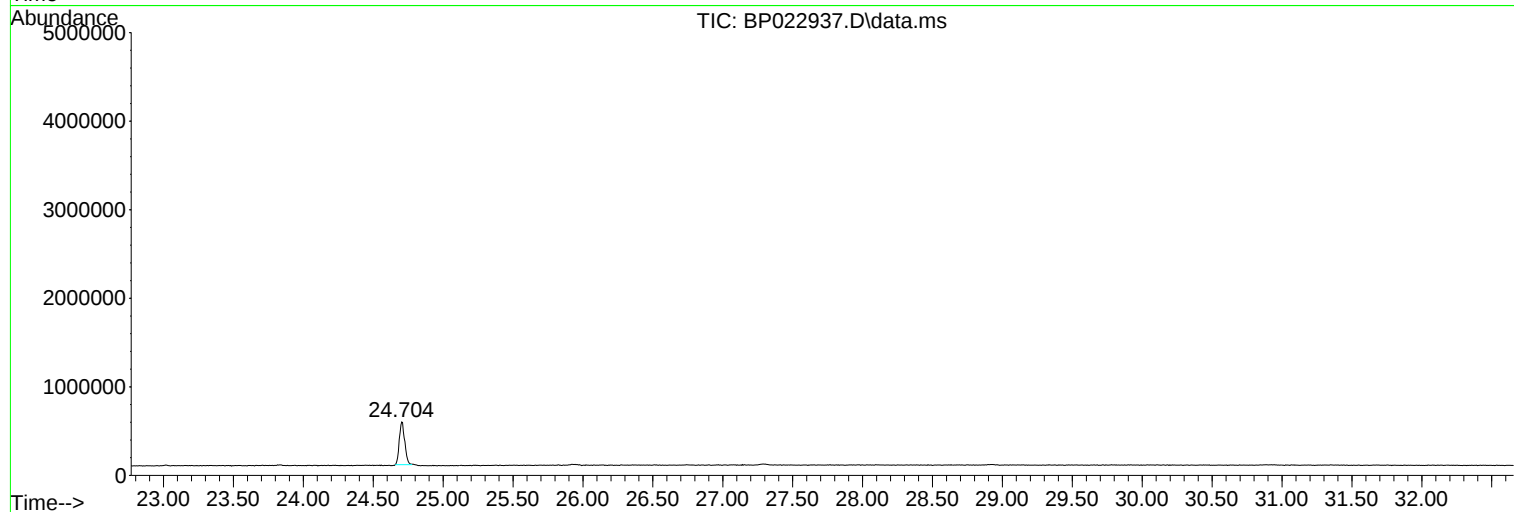
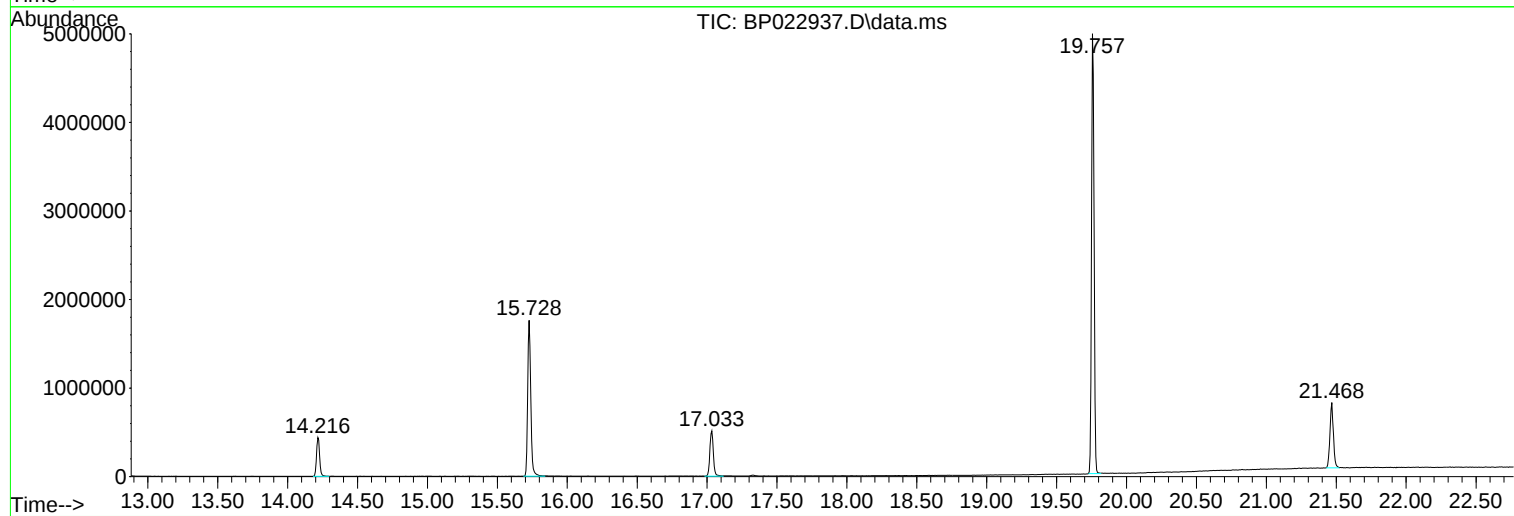
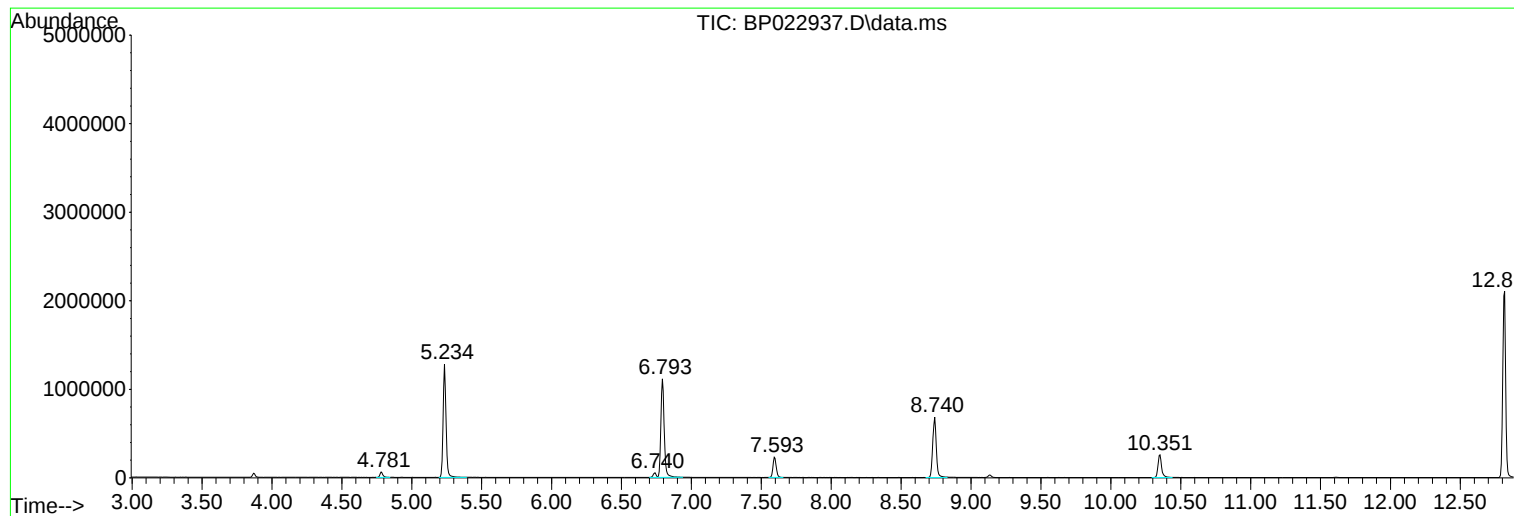
Sum of corrected areas: 21891329

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110724.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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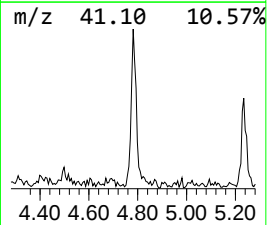
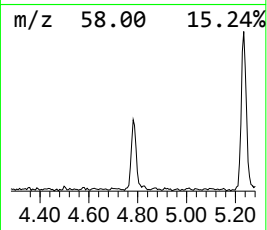
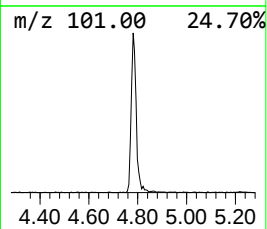
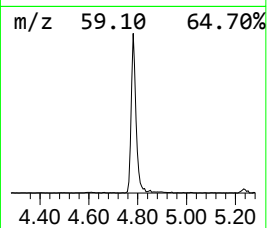
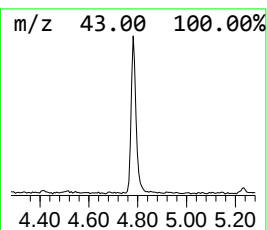
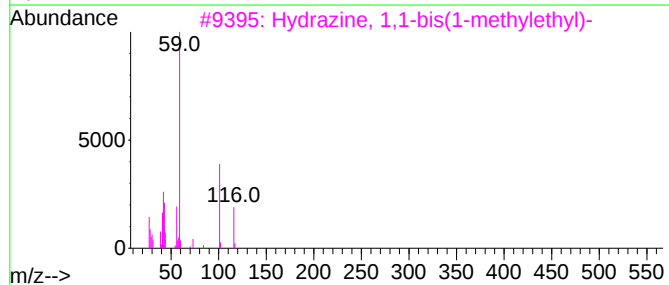
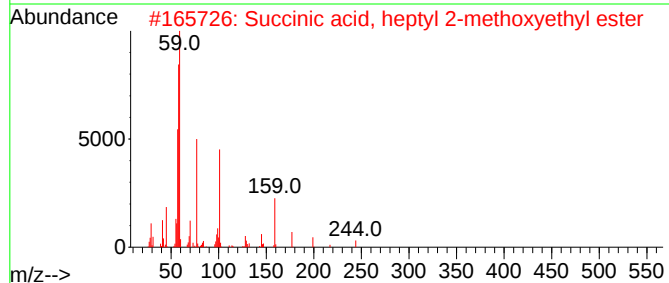
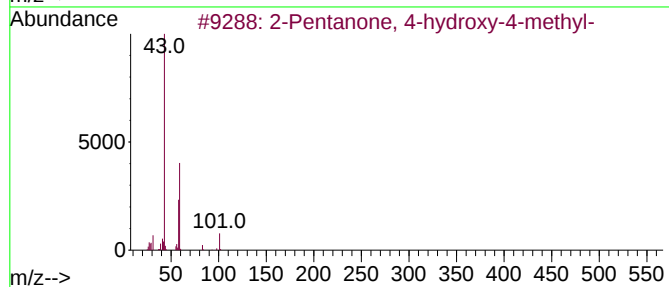
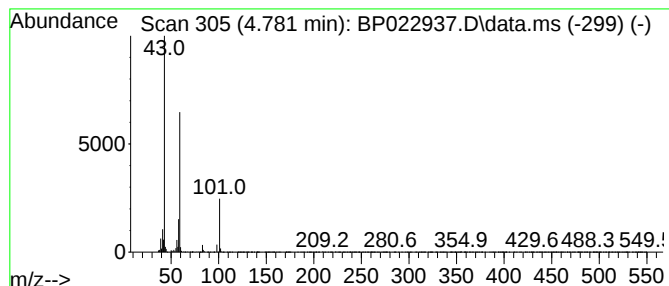
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.781	4.70 ng	87704	1,4-Dichlorobenzene-d4	7.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	40		
3	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38		
4	Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38		



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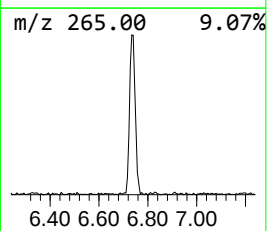
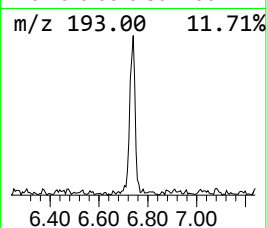
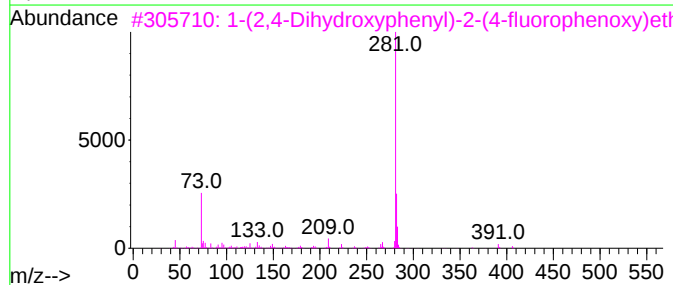
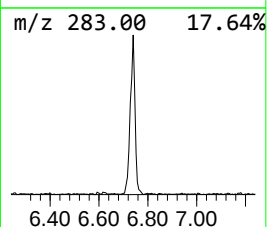
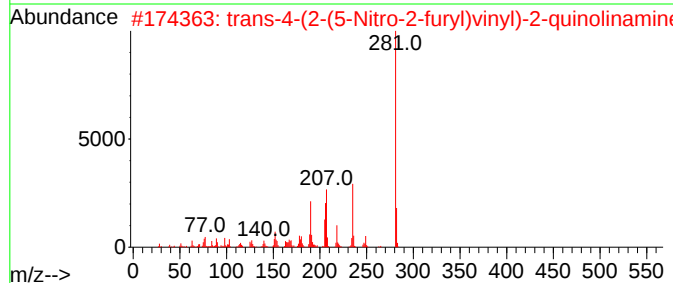
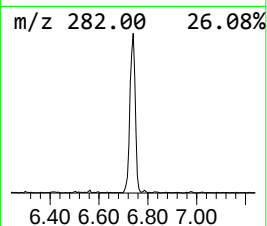
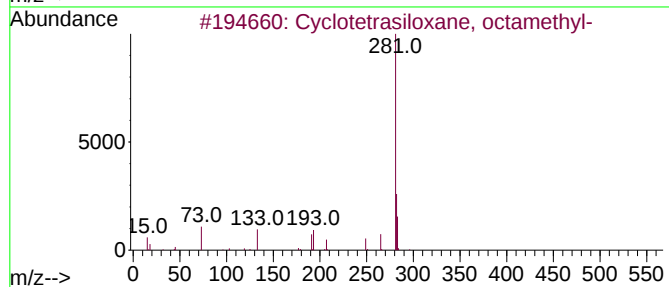
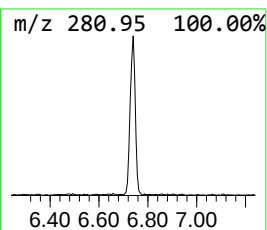
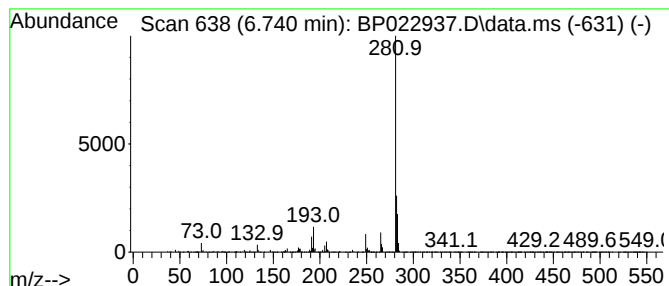
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Peak Number 2 Cyclotetrasiloxane, octamet... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.740	4.17 ng	77774	1,4-Dichlorobenzene-d4	7.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	78
2			trans-4-(2-(5-Nitro-2-furyl)vinyl)-2-(4-fluorophenoxy)ethyl	281	C15H11N3O3	000847-10-9	53
3			1-(2,4-Dihydroxyphenyl)-2-(4-fluorophenoxy)ethyl	406	C20H27F04Si2	1000474-42-3	53
4			2,4-Dihydroxy-3-methylbenzaldehyde	296	C14H24O3Si2	1000514-43-5	50
5			dibenz[b,f]azocine, 6-phenyl-	281	C21H15N	1000397-18-9	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.781	4.7	ng	87704	1	7.593	373285	20.0
Cyclotetrasilox...	6.740	4.2	ng	77774	1	7.593	373285	20.0