Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 25 00:20:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Dec 24 00:19:17 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
DBYC0

Manual IntegrationsAPPROVED

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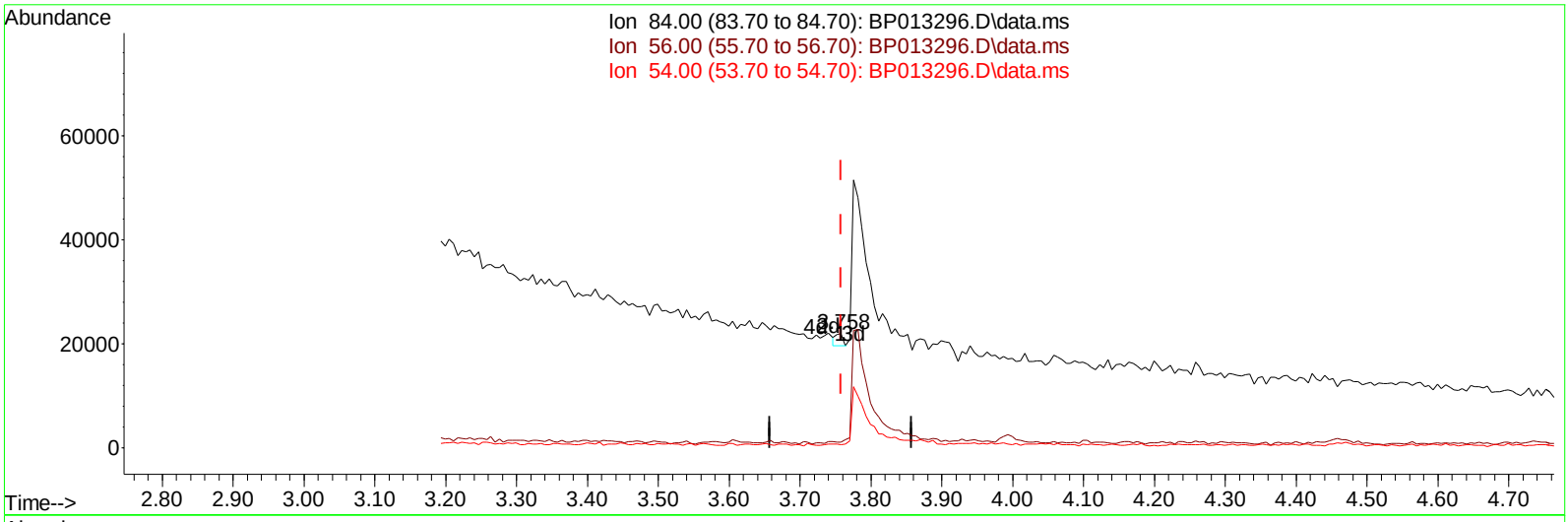
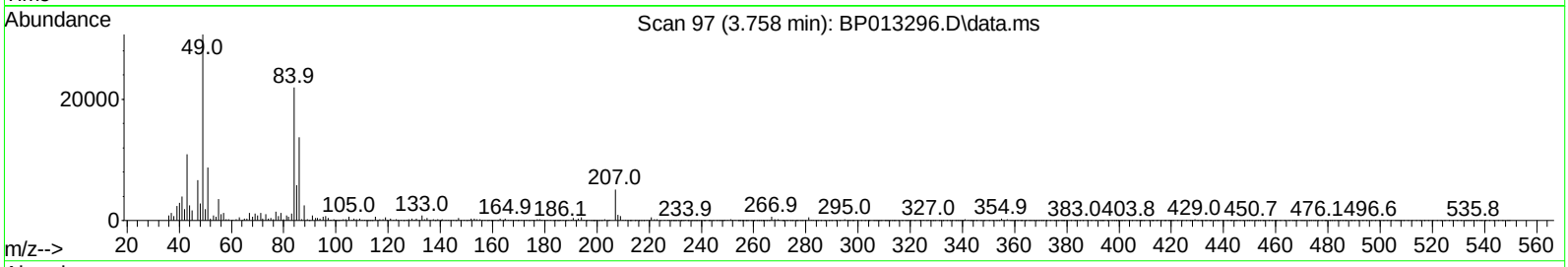
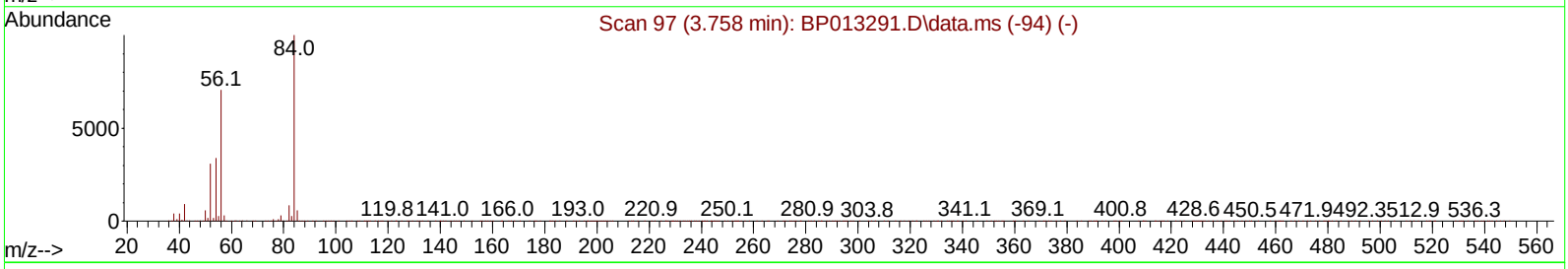
Supervised By :Yogesh Patel 12/25/2022







TIC: BP013296.D\data.ms

(4) Pyridine-d5 (S)

3.758min (+ 0.000) 0.03 ng/ul

response 1557

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	4.63#
54.00	30.50	2.70#
0.00	0.00	0.00

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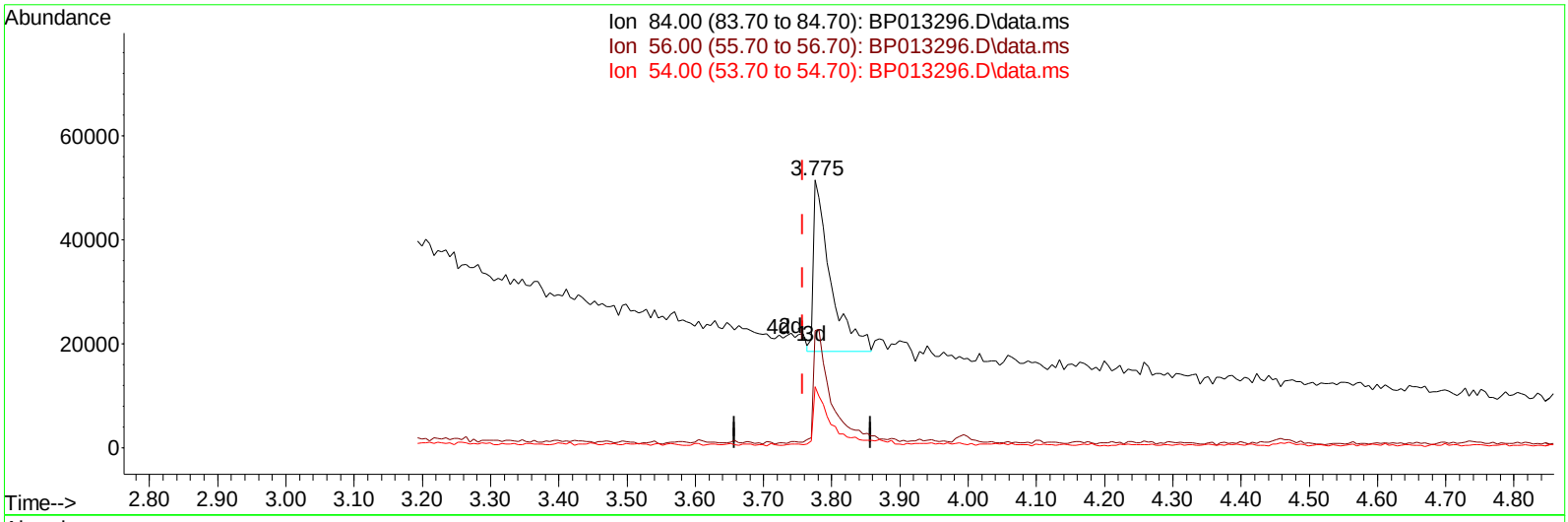

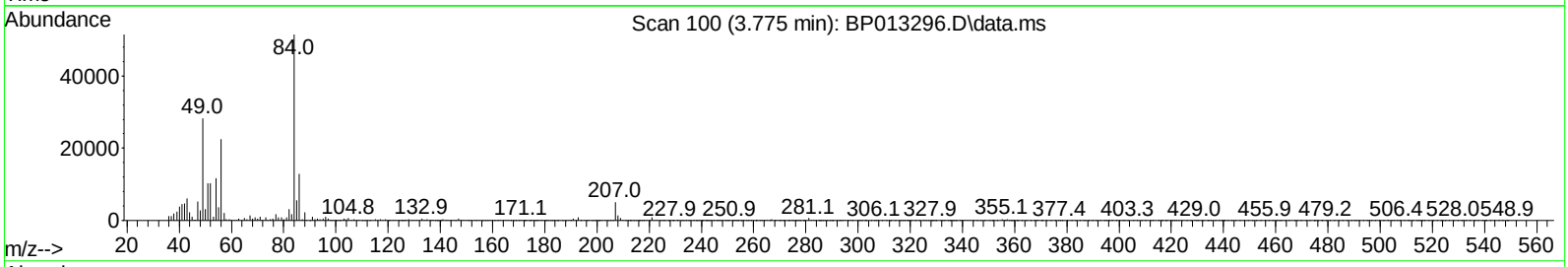

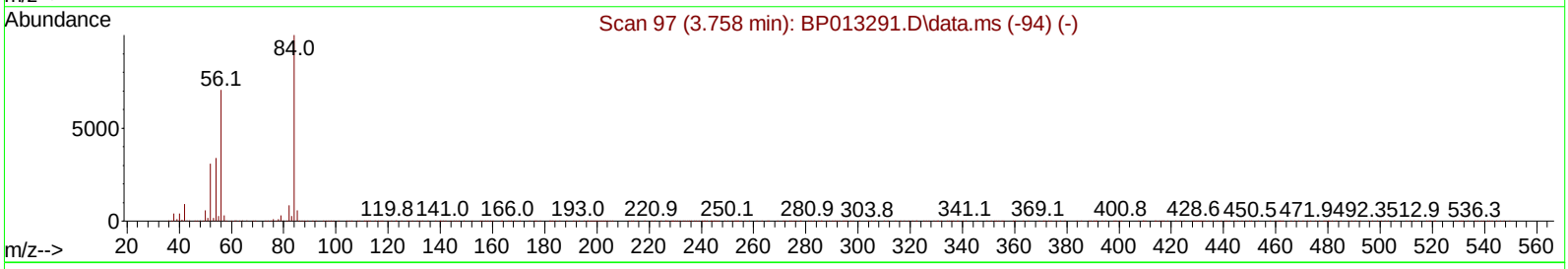

TIC: BP013296.D\data.ms

(4) Pyridine-d5 (S)

3.775min (+ 0.018) 1.26 ng/ul m

response 57639

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	43.74#
54.00	30.50	22.75#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	689289	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	2972656	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	1791186	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	3192139	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2699511	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	3229058	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	49134	3.048	ng/uL	0.00
4) Pyridine-d5	3.775	84	57639m	1.259	ng/ul	0.02
7) Phenol-d5	7.105	99	984218	17.530	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	640737	18.918	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	825034	18.667	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	654546	14.228	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	411170	18.826	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	461198	18.757	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	830232	17.383	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	758420	11.248	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	2500890	18.167	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	2764397	18.275	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	221903	8.961	ng/ul	0.00
60) Fluorene-d10	15.581	176	2054348	17.640	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	182521	8.885	ng/ul	0.00
73) Anthracene-d10	17.445	188	2564091	17.805	ng/ul	0.00
81) Pyrene-d10	19.675	212	2915303	18.814	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	3065909	19.049	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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