Quantitation Report (Qedit)

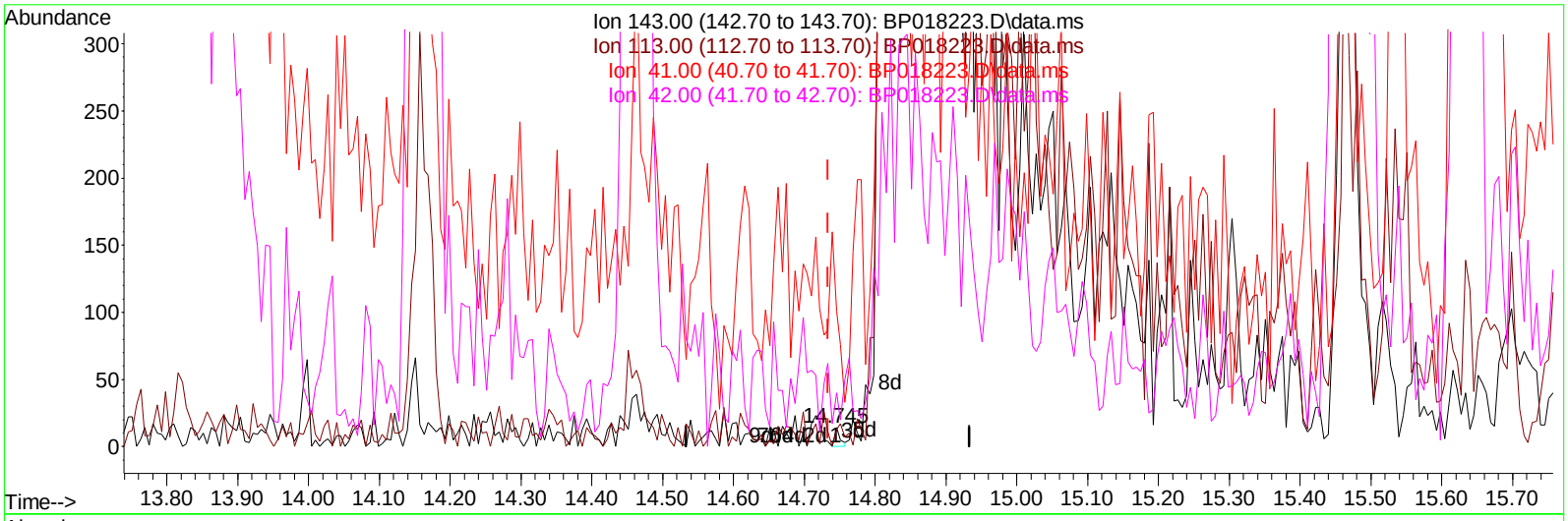
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018223.D
Acq On : 17 Nov 2023 16:12
Operator : MA/JU
Sample : 05369-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

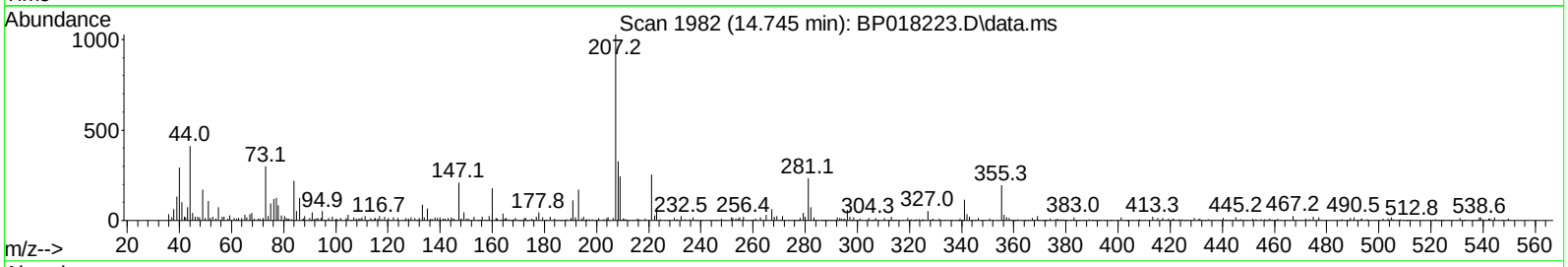
Quant Time: Nov 17 21:45:30 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 18:13:37 2023
Response via : Initial Calibration

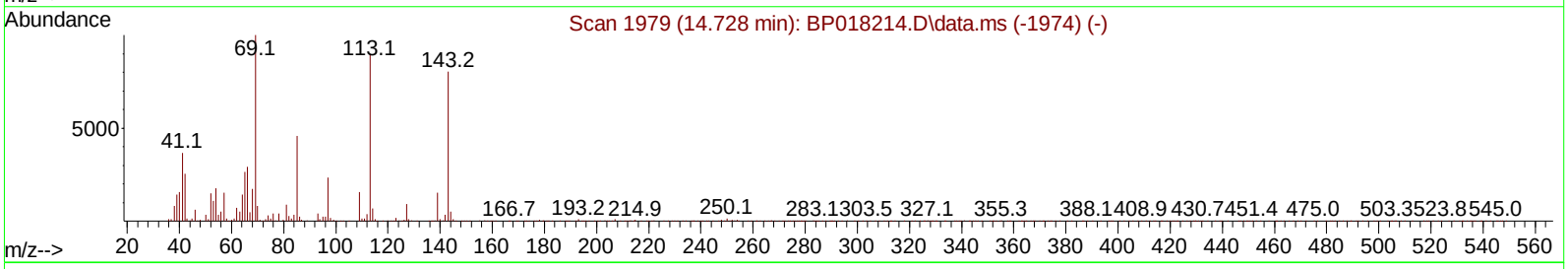
Instrument :
BNA_P
ClientSampleId :
GCDN3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2023
Supervised By :mohammad ahmed 11/19/2023







TIC: BP018223.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.012) 0.00 ng/ul

response 8

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	92.86
41.00	47.70	721.43#
42.00	30.10	285.71#

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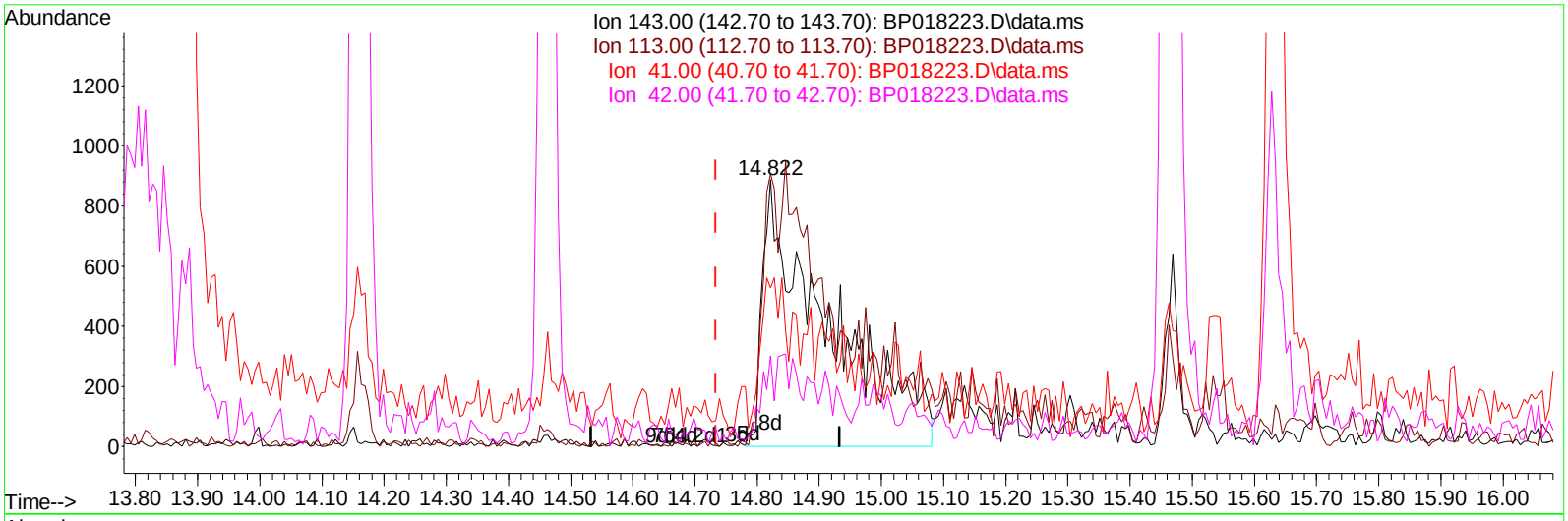

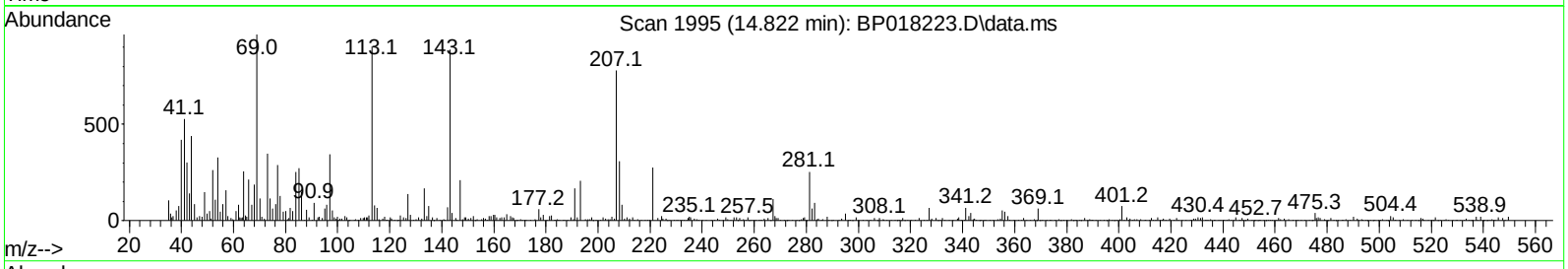

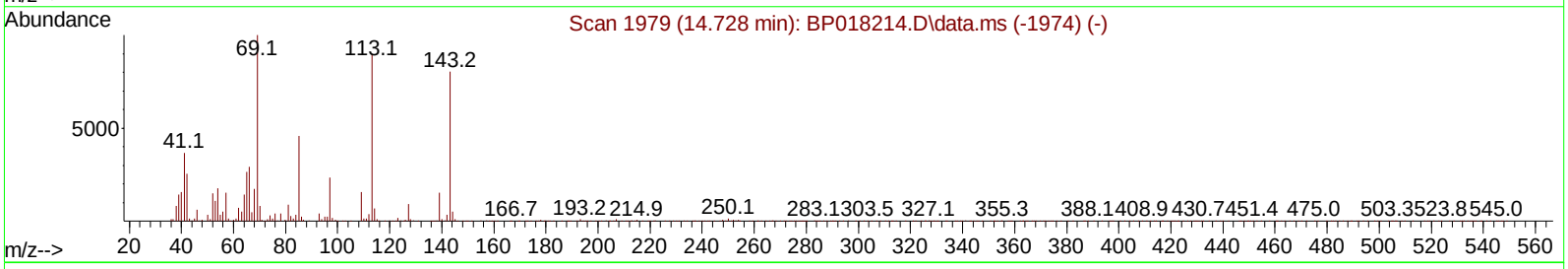

TIC: BP018223.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.088) 3.91 ng/ul m

response 6491

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	102.14
41.00	47.70	59.53#
42.00	30.10	34.16

Quantitation Report (QT Reviewed)

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Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	53668	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	204719	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	120452	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	269252	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	269629	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	310826	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	5790	3.806	ng/uL	0.00
4) Pyridine-d5	3.599	84	39194	9.881	ng/ul	0.00
7) Phenol-d5	6.969	99	26909	5.194	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.134	67	109633	36.006	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	103146	25.139	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	61886	14.576	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	62738	34.015	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	60480	28.944	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	97871	26.280	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	144150	28.225	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	414844	37.327	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	419175	34.628	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	6491m	3.909	ng/ul	0.09
60) Fluorene-d10	15.463	176	312942	34.105	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	54506	28.002	ng/ul	0.00
73) Anthracene-d10	17.345	188	573926	39.412	ng/ul	0.00
81) Pyrene-d10	19.604	212	726455	40.336	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	748313	41.404	ng/ul	0.00

Target Compounds Qvalue

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

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