Quantitation Report (Qedit)

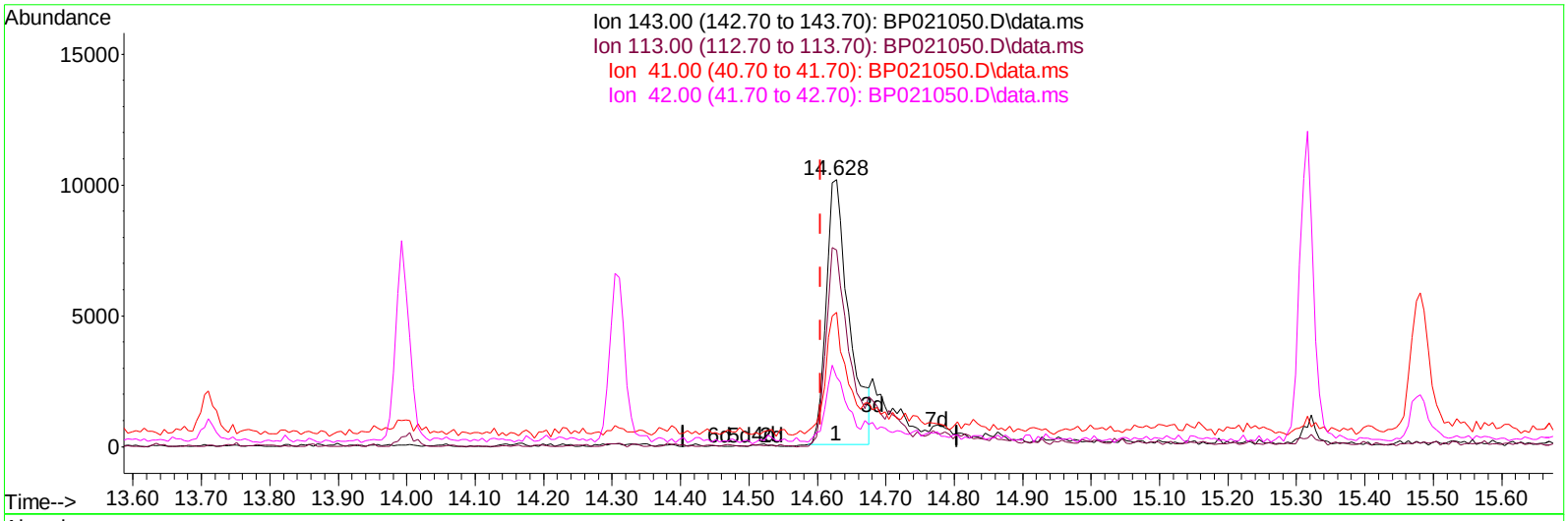
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021050.D
Acq On : 18 Jul 2024 17:36
Operator : MA/JU
Sample : P3260-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

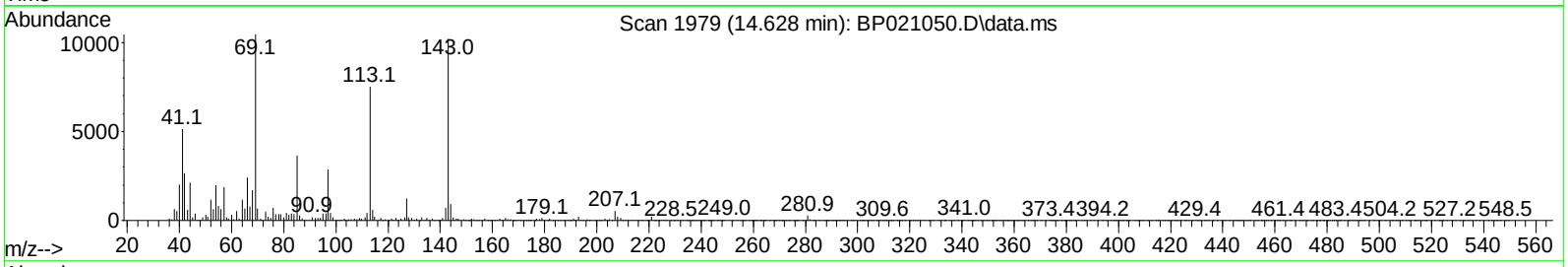
Quant Time: Jul 18 18:51:18 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 18 14:23:21 2024
Response via : Initial Calibration

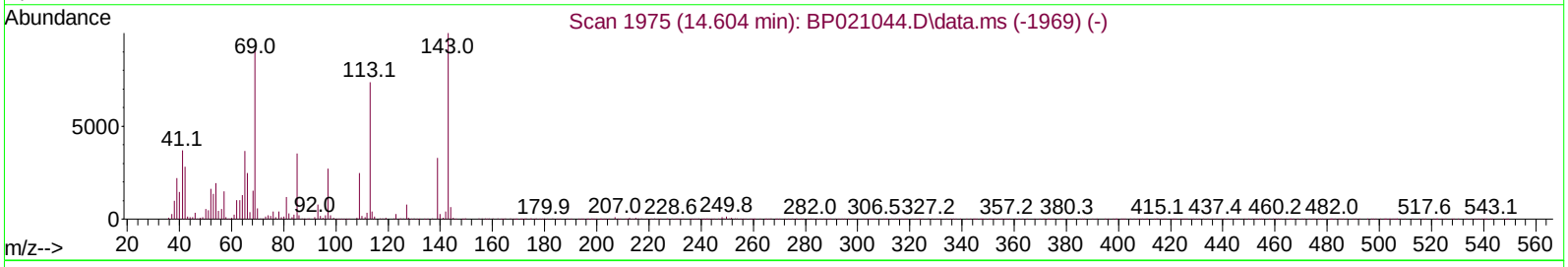
Instrument :
BNA_P
ClientSampleId :
E08F4

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/19/2024
Supervised By :mohammad ahmed 07/20/2024







TIC: BP021050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.024) 6.60 ng/ul

response 23521

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	73.69
41.00	37.90	50.37#
42.00	27.10	26.07

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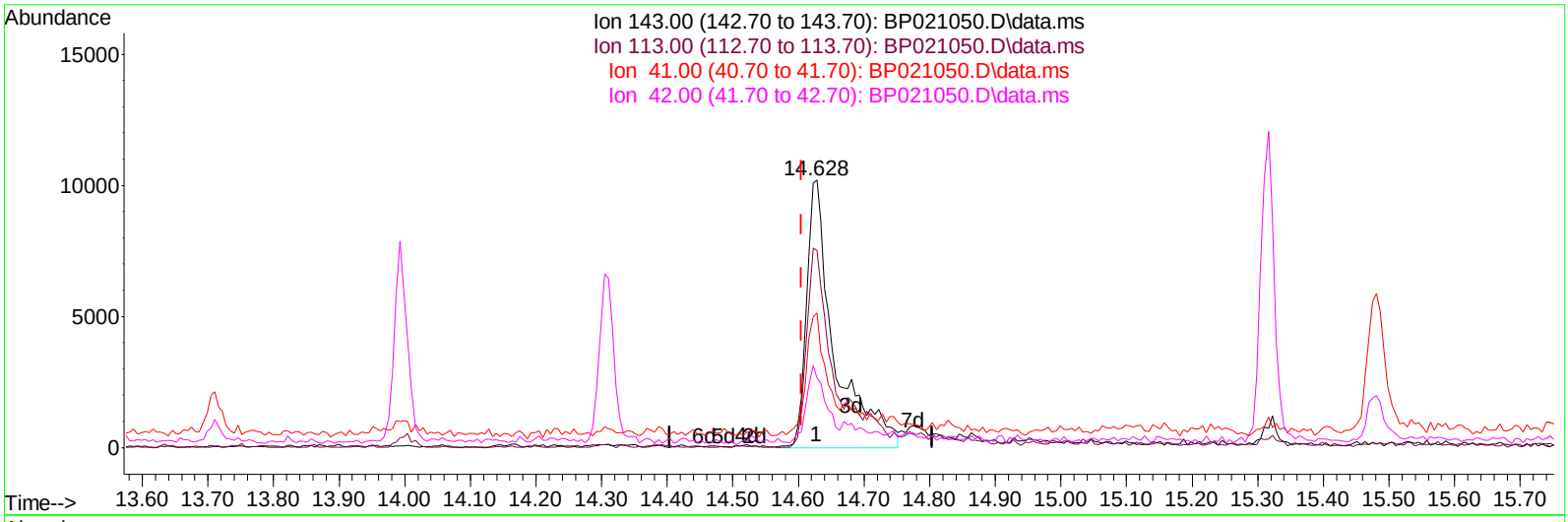

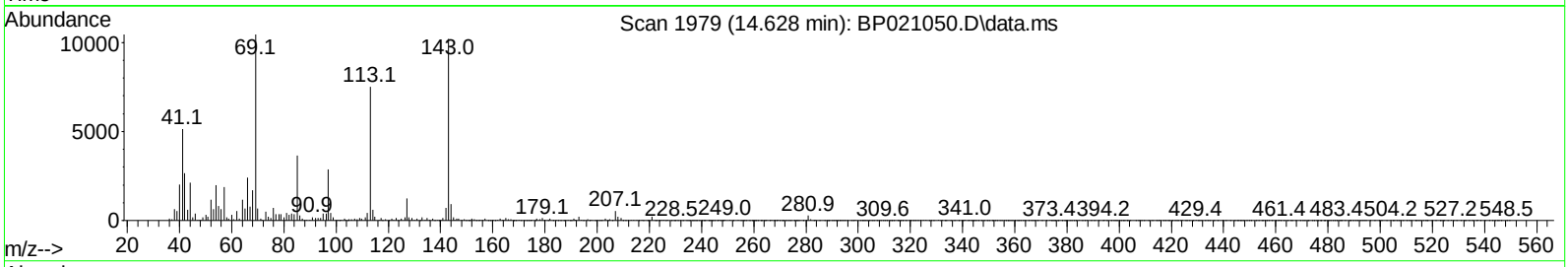

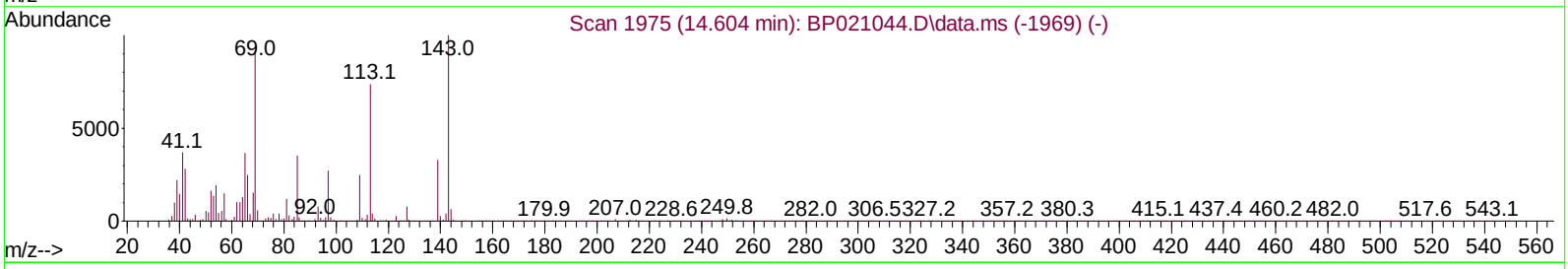

TIC: BP021050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.024) 8.46 ng/ul m

response 30160

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	73.69
41.00	37.90	50.37#
42.00	27.10	26.07

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
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Acq On : 18 Jul 2024 17:36
Operator : MA/JU
Sample : P3260-01
Misc :
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Quant Time: Jul 18 18:52:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 18 14:23:21 2024
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Instrument :
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ClientSampleId :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	140625	20.000	ng/ul	0.00
20) Naphthalene-d8	10.452	136	548271	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	344754	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.122	188	752823	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	805505	20.000	ng/ul	-0.01
88) Perylene-d12	24.892	264	1046953	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	12892	4.172	ng/uL	0.00
4) Pyridine-d5	3.658	84	47719	5.334	ng/ul	0.00
7) Phenol-d5	6.905	99	87760	7.655	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	200058	28.794	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	253379	26.824	ng/ul	0.00
15) 4-Methylphenol-d8	8.411	113	163399	17.160	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	137381	32.260	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	156720	32.154	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.105	165	262738	29.811	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	247231	21.175	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	804890	32.116	ng/ul	0.00
49) Acenaphthylene-d8	13.993	160	934870	33.221	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	30160m	8.458	ng/ul	0.02
60) Fluorene-d10	15.316	176	716727	34.859	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	126744	27.825	ng/ul	0.00
73) Anthracene-d10	17.222	188	1241882	37.141	ng/ul	0.00
81) Pyrene-d10	19.616	212	1596840	32.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.651	264	1961125	37.980	ng/ul	-0.04

Target Compounds	Qvalue

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

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