

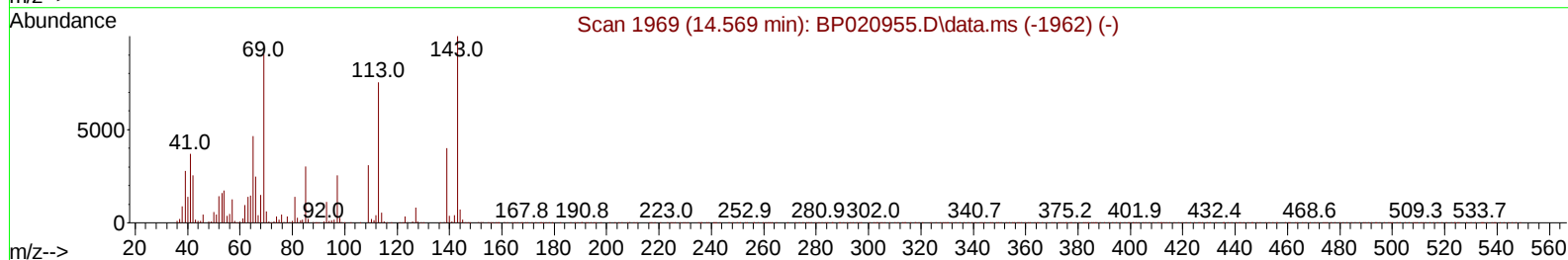
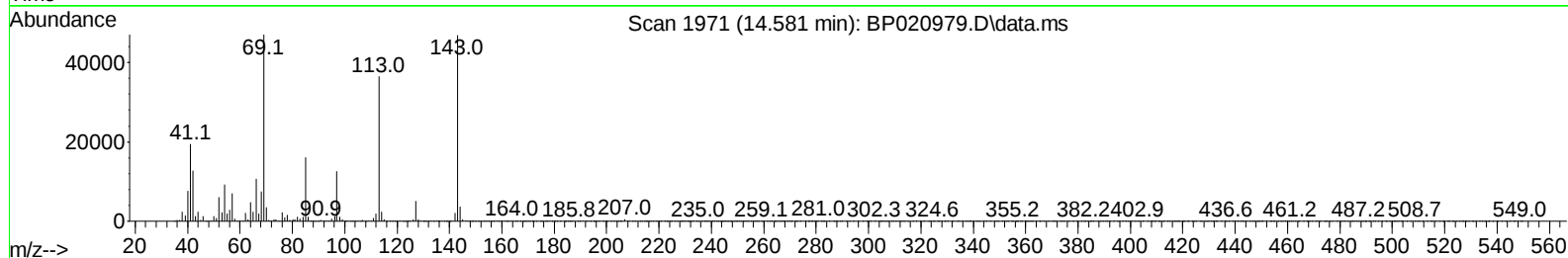
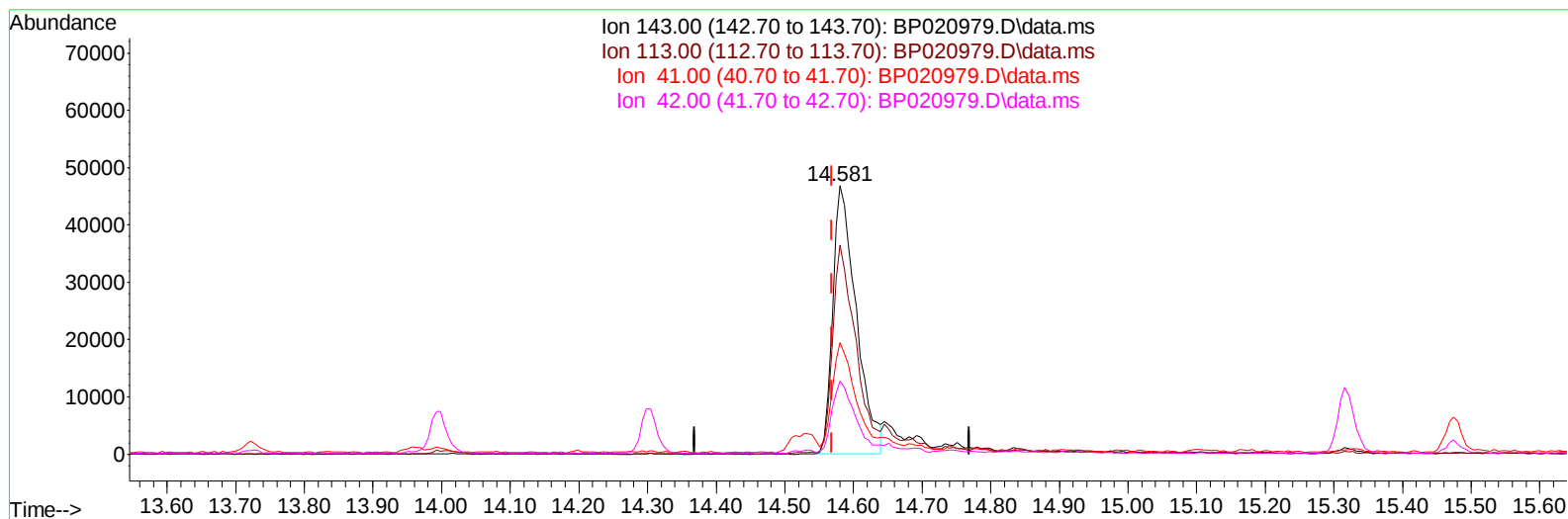
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020979.D
Acq On : 16 Jul 2024 06:27
Operator : MA/JU
Sample : P3100-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BW4

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/16/2024
Supervised By :mohammad ahmed 07/18/2024

Quant Time: Jul 16 06:53:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 15 12:22:22 2024
Response via : Initial Calibration



TIC: BP020979.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.012) 24.09 ng/u1

response 111043

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	77.83
41.00	37.90	41.70
42.00	27.10	27.37

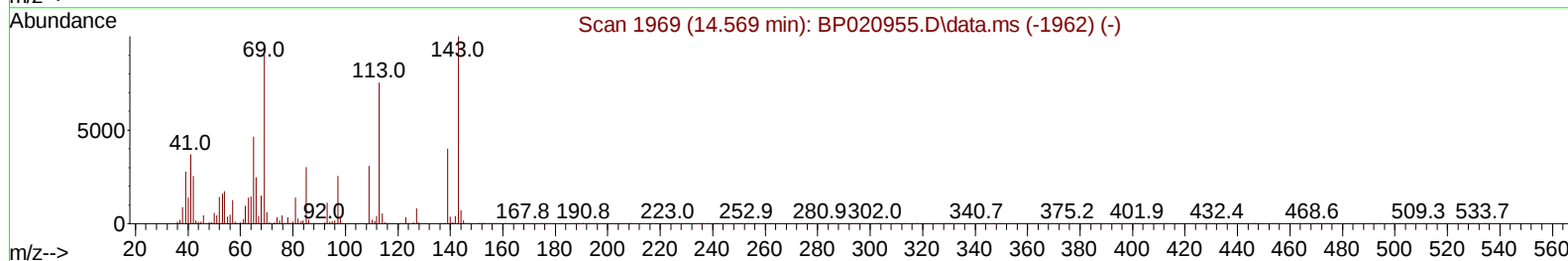
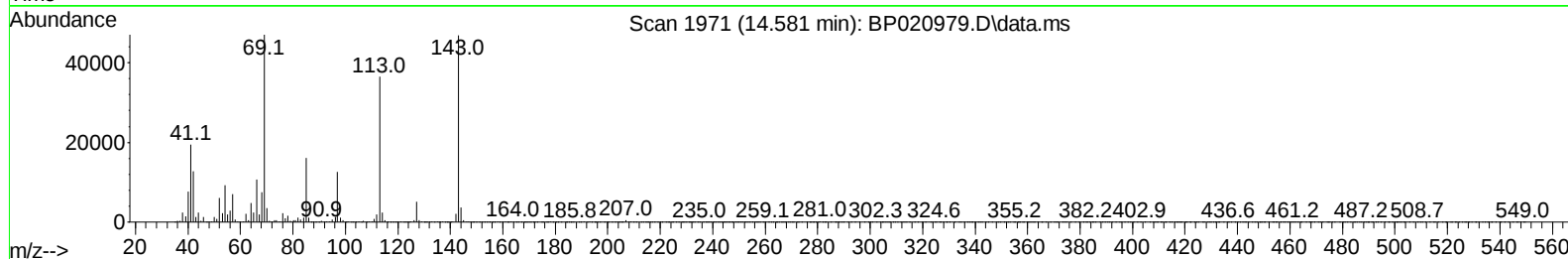
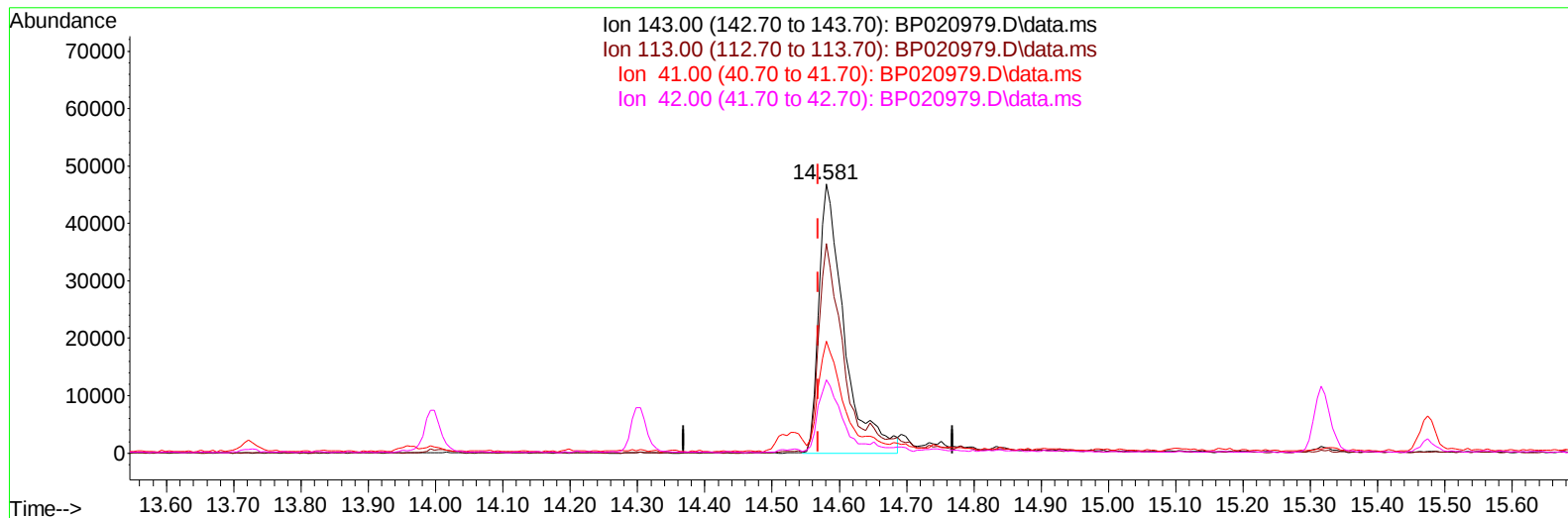
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TIC: BP020979.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.012) 26.47 ng/ul m

response 122045

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	77.83
41.00	37.90	41.70
42.00	27.10	27.37

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 Sample : P3100-07
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 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/16/2024
 Supervised By :mohammad ahmed 07/18/2024

Quant Time: Jul 16 06:54:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 15 12:22:22 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	191235	20.000	ng/ul	0.01
20) Naphthalene-d8	10.451	136	746517	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.304	164	445756	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.122	188	917036	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	999562	20.000	ng/ul	0.00
88) Perylene-d12	24.839	264	1171979	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	18399	4.378	ng/uL	0.00
4) Pyridine-d5	3.664	84	269368	22.140	ng/ul	0.01
7) Phenol-d5	6.905	99	410166	26.308	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.046	67	249834	26.442	ng/ul	0.01
11) 2-Chlorophenol-d4	7.246	132	344491	26.818	ng/ul	0.01
15) 4-Methylphenol-d8	8.410	113	336683	26.001	ng/ul	0.02
21) Nitrobenzene-d5	8.857	128	162984	28.109	ng/ul	0.03
24) 2-Nitrophenol-d4	9.557	143	191278	28.822	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.098	165	320619	26.718	ng/ul	0.01
31) 4-Chloroaniline-d4	10.604	131	363336	22.855	ng/ul	0.02
46) Dimethylphthalate-d6	13.722	166	972360	30.007	ng/ul	0.01
49) Acenaphthylene-d8	13.998	160	1105645	30.387	ng/ul	0.00
54) 4-Nitrophenol-d4	14.581	143	122045m	26.472	ng/ul	0.01
60) Fluorene-d10	15.316	176	819698	30.834	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	120812	21.774	ng/ul	0.01
73) Anthracene-d10	17.216	188	1249482	30.677	ng/ul	0.00
81) Pyrene-d10	19.604	212	1659780	26.859	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.621	264	1795980	31.071	ng/ul	0.02

Target Compounds Qvalue

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

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