

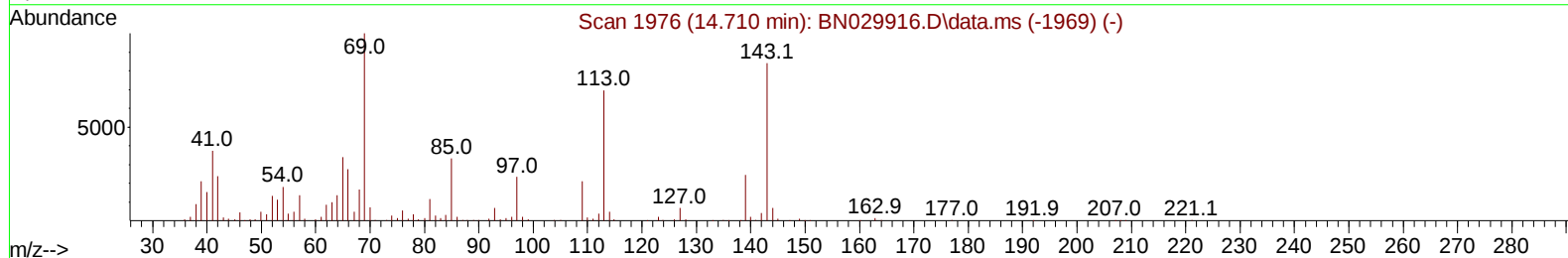
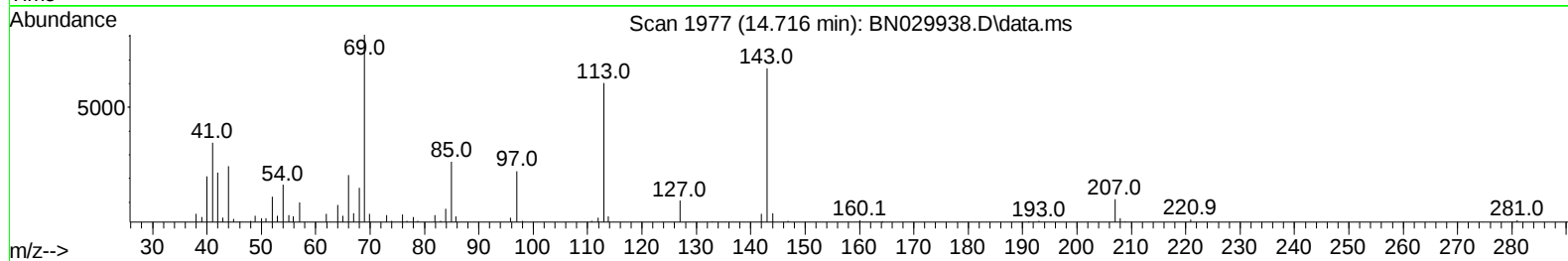
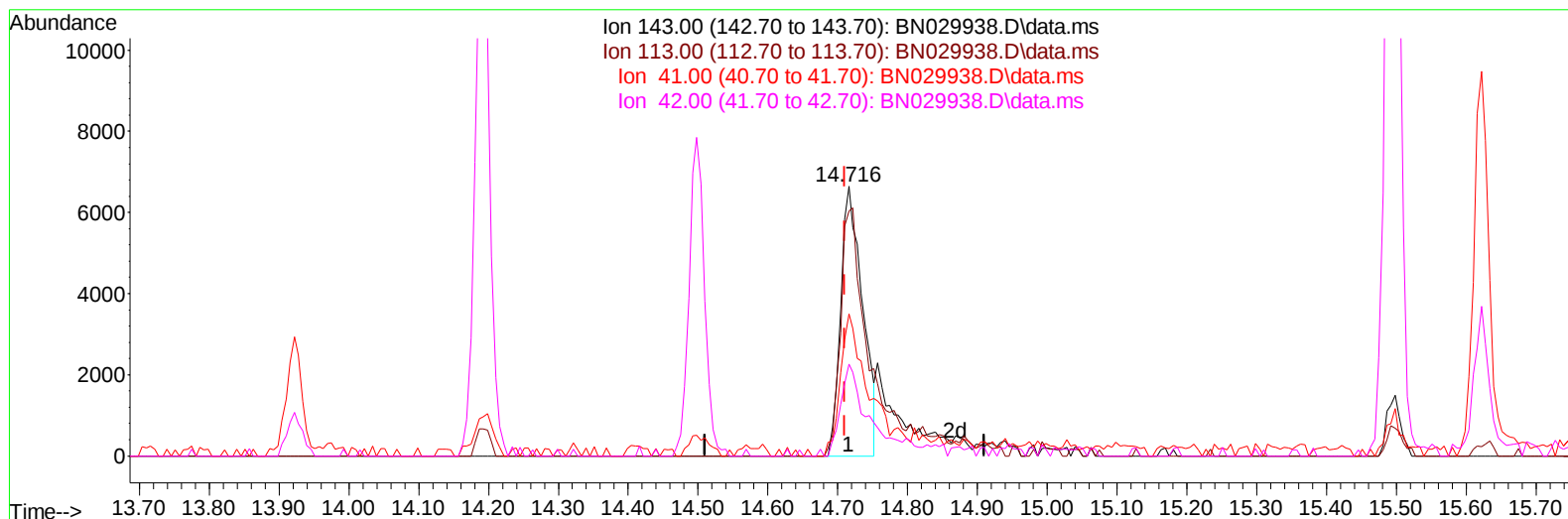
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022124\
Data File : BN029938.D
Acq On : 22 Feb 2024 21:38
Operator : MA/JU
Sample : P1494-05
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH3Y5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/23/2024
Supervised By :mohammad ahmed 02/24/2024

Quant Time: Feb 22 22:44:00 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 21 22:34:54 2024
Response via : Initial Calibration



TIC: BN029938.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 3.37 ng/u1

response 14402

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	90.95
41.00	45.90	52.77
42.00	28.40	34.02

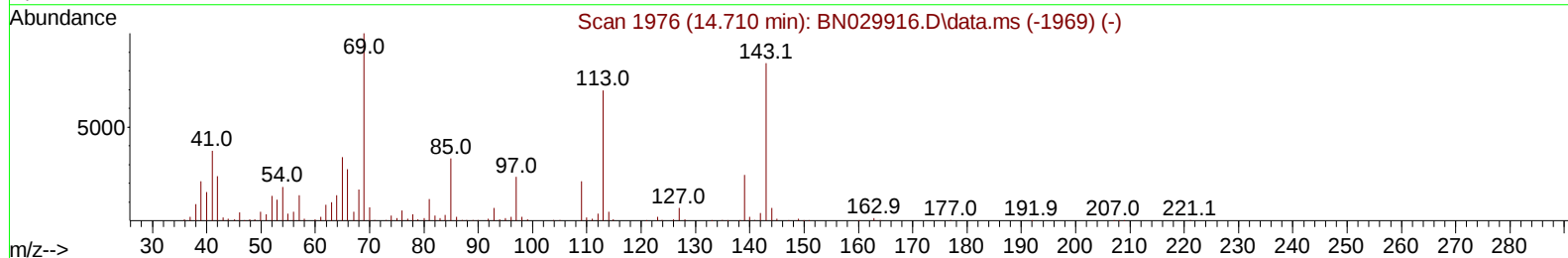
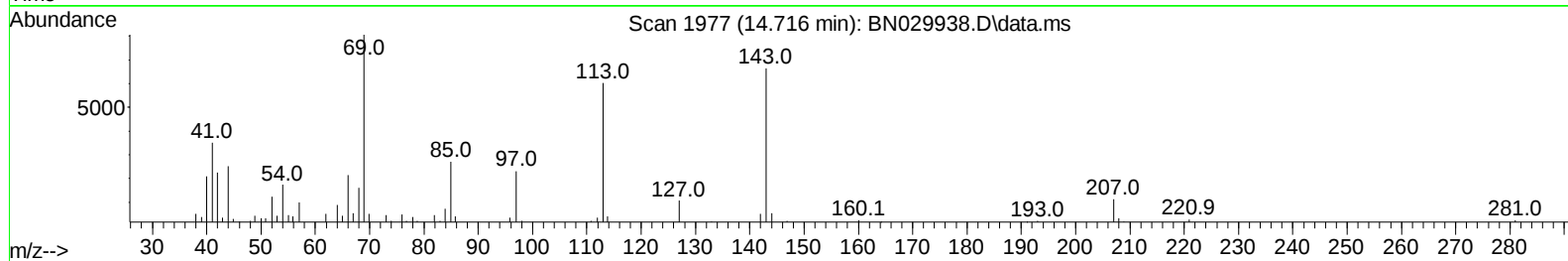
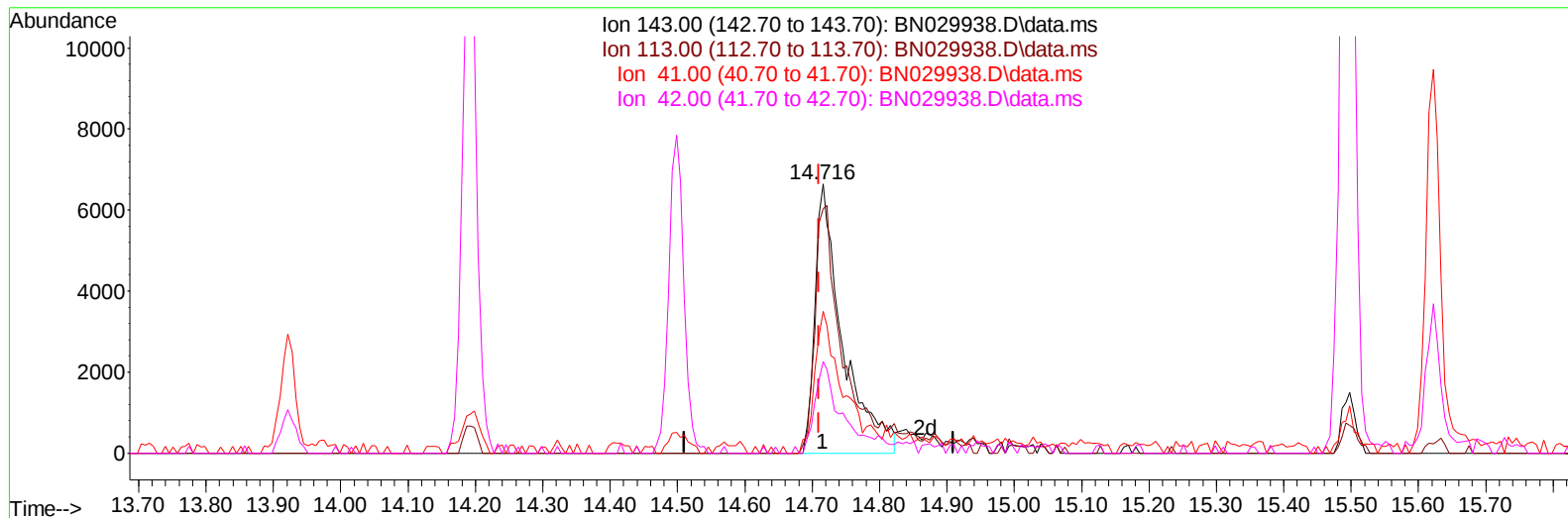
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TIC: BN029938.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 4.40 ng/ul m

response 18834

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	90.95
41.00	45.90	52.77
42.00	28.40	34.02

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/23/2024
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Quant Time: Feb 22 23:00:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 21 22:34:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	137186	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	568890	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	302659	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	644118	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	520331	20.000	ng/ul	0.00
88) Perylene-d12	23.815	264	600778	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	16540	3.966	ng/uL	0.00
4) Pyridine-d5	3.734	84	117794	10.590	ng/ul	0.00
7) Phenol-d5	7.022	99	102201	7.965	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	351968	41.100	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	310805	33.425	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	188972	19.747	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	199069	45.086	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	186913	48.607	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	304917	40.172	ng/ul	0.00
31) 4-Chloroaniline-d4	10.804	131	469605	36.266	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	1043469	46.308	ng/ul	0.00
49) Acenaphthylene-d8	14.192	160	1222221	45.674	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	18834m	4.404	ng/ul	0.00
60) Fluorene-d10	15.498	176	927669	48.393	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	117118	37.316	ng/ul	0.00
73) Anthracene-d10	17.351	188	1538475	50.965	ng/ul	0.00
81) Pyrene-d10	19.651	212	1698279	54.388	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.662	264	1553439	51.674	ng/ul	0.00

Target Compounds Qvalue

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

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