

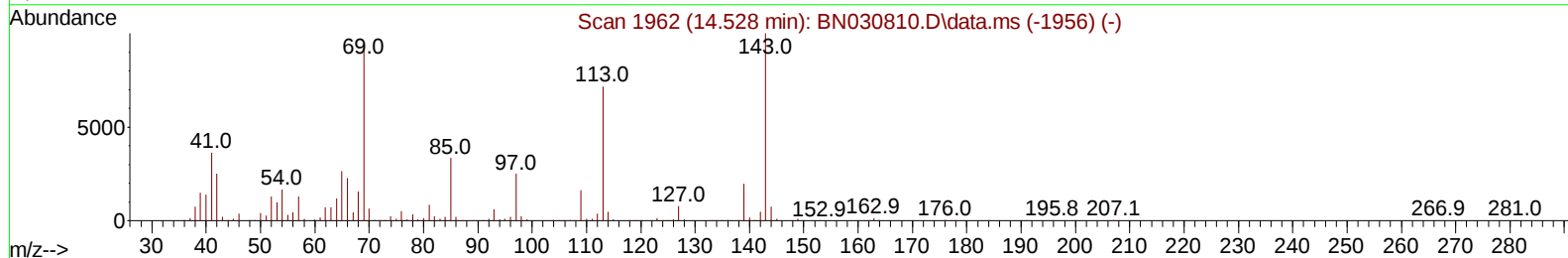
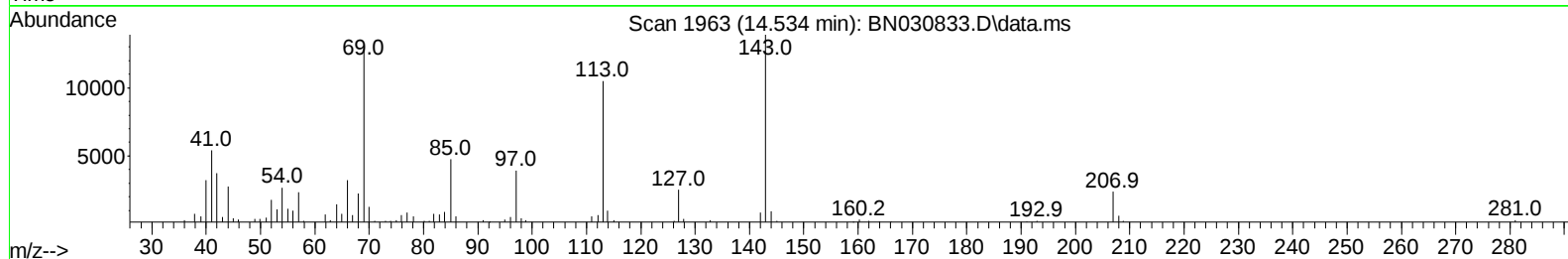
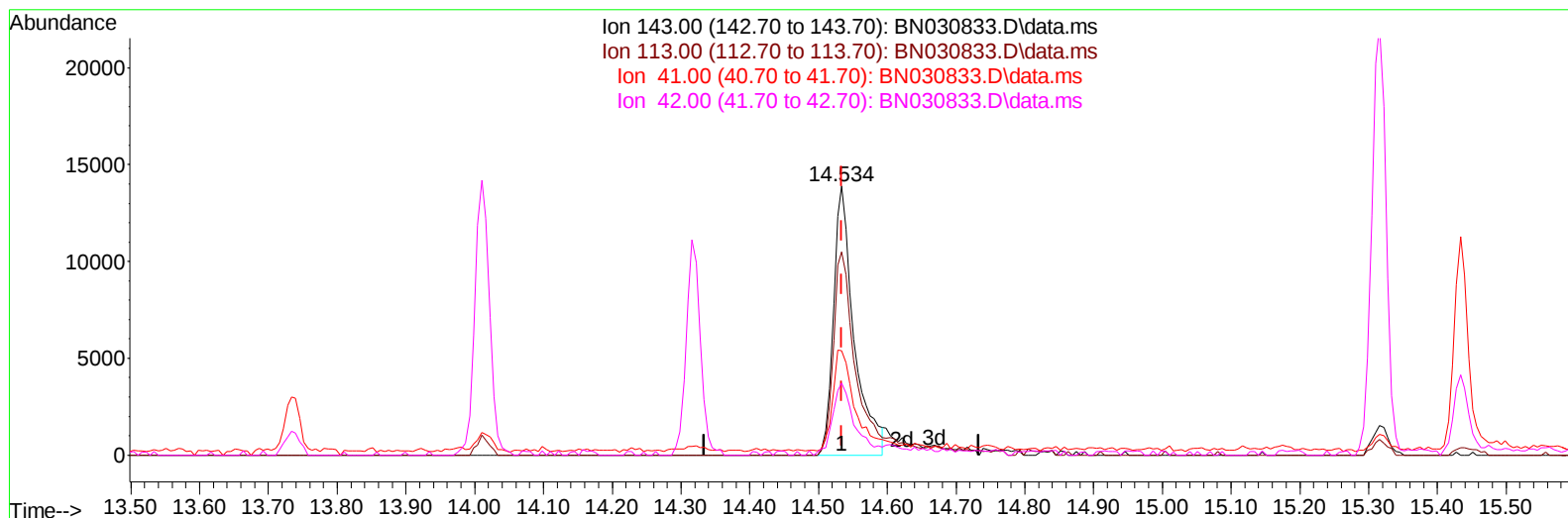
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041024\
 Data File : BN030833.D
 Acq On : 11 Apr 2024 02:24
 Operator : MA/JU
 Sample : P2035-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AN7

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:57:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 09 12:10:11 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/12/2024
 Supervised By :mohammad ahmed 04/12/2024



TIC: BN030833.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.534min (+ 0.000) 4.11 ng/u1

response 28896

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	72.10	75.55
41.00	38.20	38.87
42.00	26.70	26.88

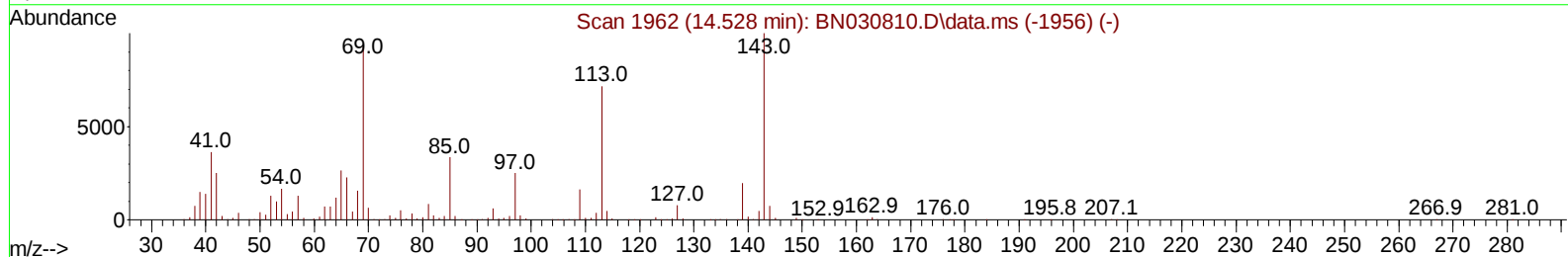
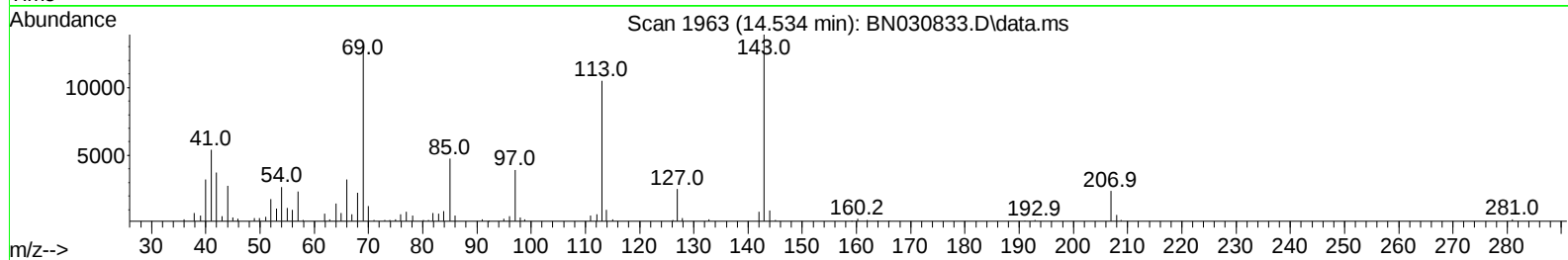
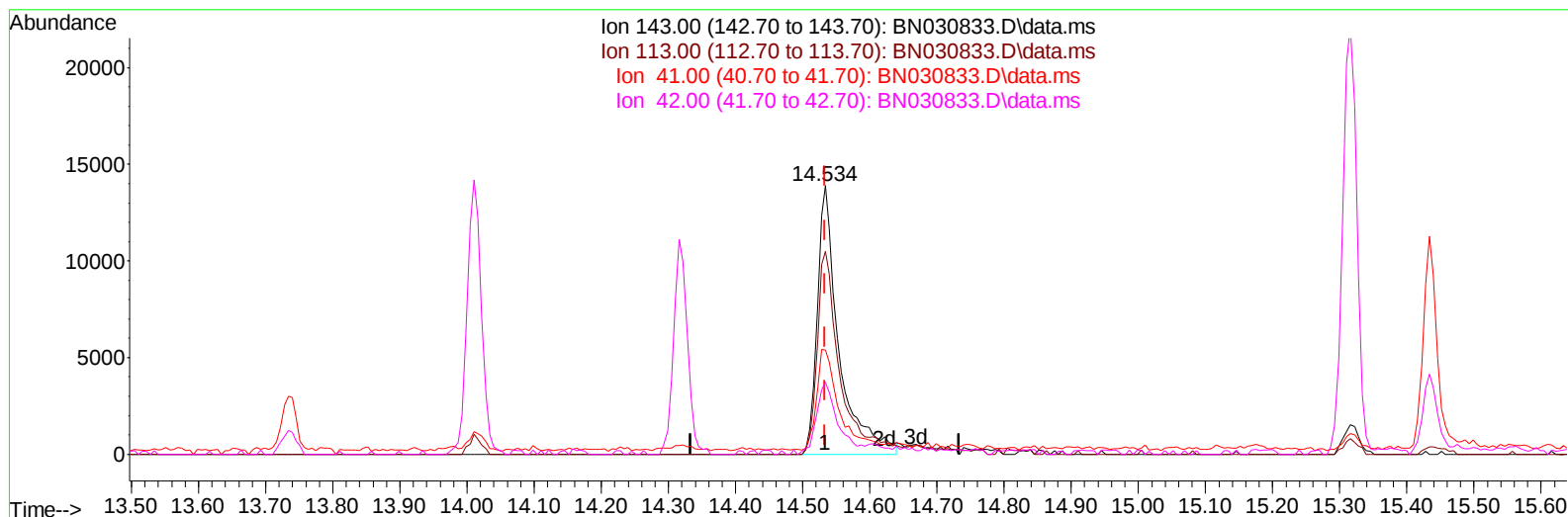
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TIC: BN030833.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.534min (+ 0.000) 4.44 ng/ul m

response 31205

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	72.10	75.55
41.00	38.20	38.87
42.00	26.70	26.88

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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	173975	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	798523	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	549971	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	1272200	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1333104	20.000	ng/ul	0.00
88) Perylene-d12	24.233	264	1408014	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	15762	4.248	ng/uL	0.00
4) Pyridine-d5	3.599	84	124349	11.177	ng/ul	0.00
7) Phenol-d5	6.846	99	96989	6.891	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	321977	37.040	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	314056	29.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	197655	16.757	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	225324	40.677	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	212872	42.008	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	399315	35.994	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	537237	29.415	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	1522170	41.454	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160	1770438	41.935	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	31205m	4.443	ng/ul	0.00
60) Fluorene-d10	15.316	176	1370344	42.647	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.434	200	173131	30.831	ng/ul	0.00
73) Anthracene-d10	17.169	188	2304233	42.037	ng/ul	0.00
81) Pyrene-d10	19.469	212	2873183	40.087	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.039	264	2843358	43.826	ng/ul	0.00

Target Compounds Qvalue

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

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