

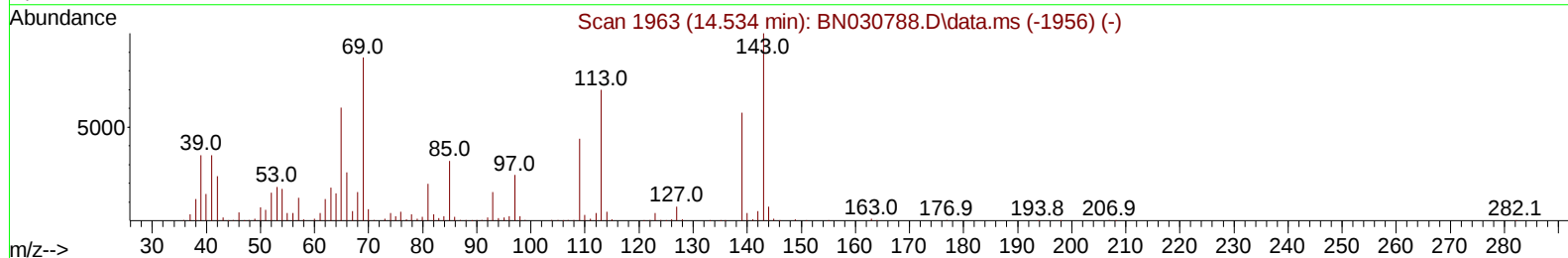
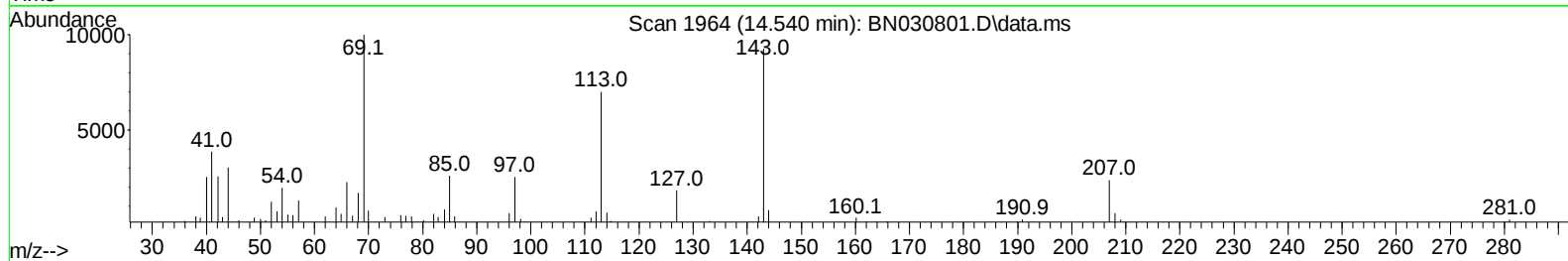
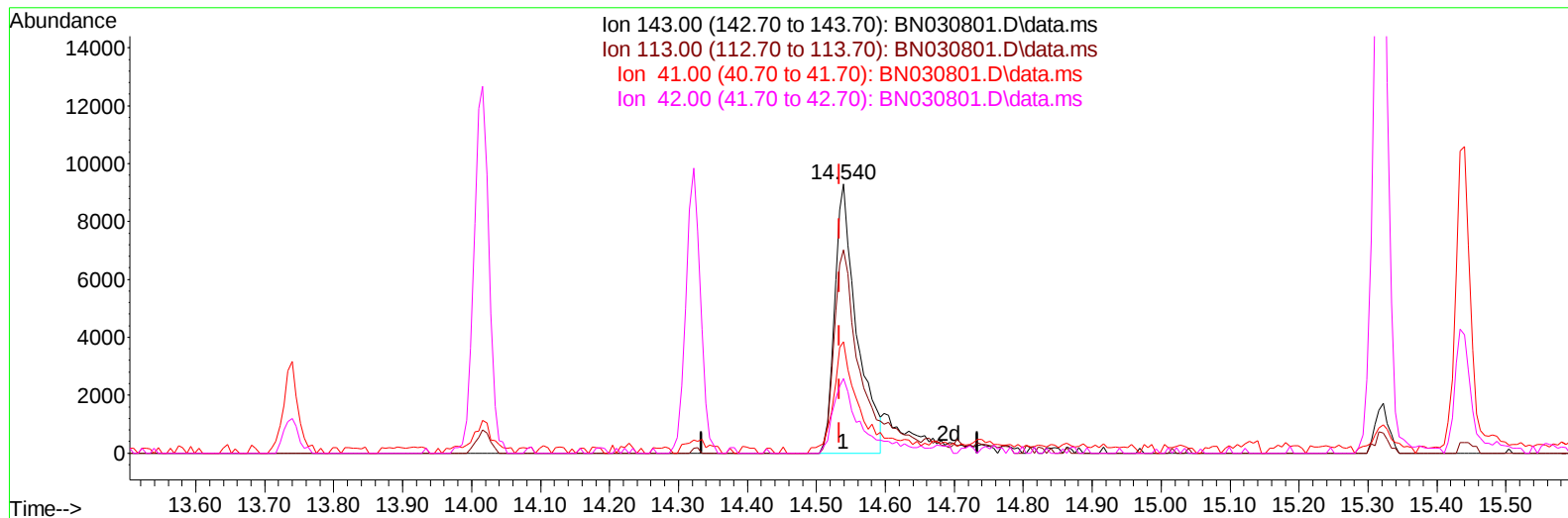
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040924\
Data File : BN030801.D
Acq On : 09 Apr 2024 22:30
Operator : MA/JU
Sample : P1973-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
YCSW1

Manual IntegrationsAPPROVED

Quant Time: Apr 09 23:02:22 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 04/10/2024
Supervised By :mohammad ahmed 04/10/2024



TIC: BN030801.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (+ 0.006) 3.19 ng/ul

response 20631

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	72.10	75.58
41.00	38.20	41.43
42.00	26.70	27.72

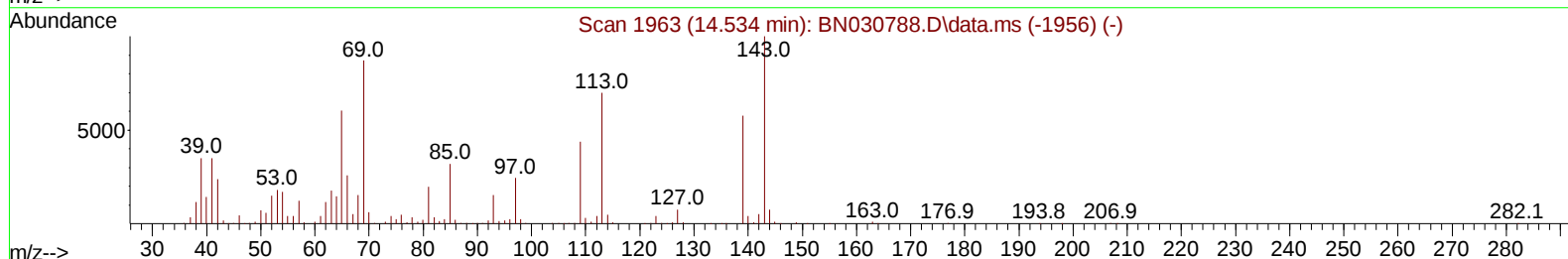
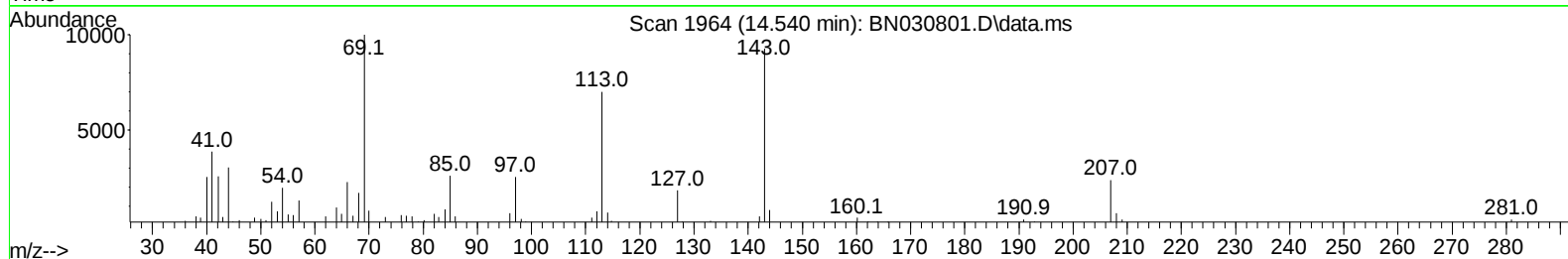
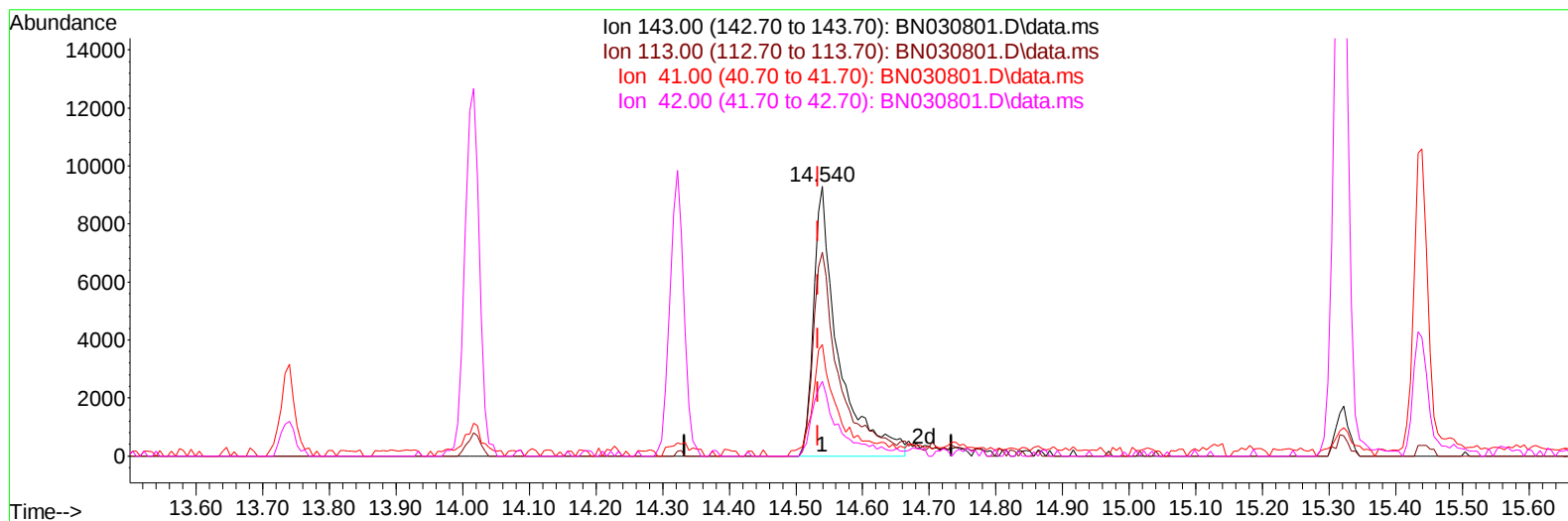
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TIC: BN030801.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (+ 0.006) 3.71 ng/ul m

response 23949

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	72.10	75.58
41.00	38.20	41.43
42.00	26.70	27.72

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 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/10/2024
 Supervised By :mohammad ahmed 04/10/2024

Quant Time: Apr 09 23:03:09 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	149838	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	705528	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.322	164	505870	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	1117848	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1077799	20.000	ng/ul	0.00
88) Perylene-d12	24.239	264	1168546	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	13443	4.206	ng/uL	0.00
4) Pyridine-d5	3.599	84	98124	10.240	ng/ul	0.00
7) Phenol-d5	6.852	99	68777	5.674	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	284631	38.019	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	263272	28.632	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	156576	15.413	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	199449	40.752	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	198443	44.322	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	344099	35.106	ng/ul	0.00
31) 4-Chloroaniline-d4	10.605	131	496795	30.786	ng/ul	0.00
46) Dimethylphthalate-d6	13.740	166	1462404	43.299	ng/ul	0.00
49) Acenaphthylene-d8	14.016	160	1682116	43.317	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	23949m	3.707	ng/ul	0.00
60) Fluorene-d10	15.322	176	1328384	44.945	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.440	200	178776	36.233	ng/ul	0.00
73) Anthracene-d10	17.169	188	2242062	46.551	ng/ul	0.00
81) Pyrene-d10	19.475	212	2663396	45.963	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.045	264	2585282	48.014	ng/ul	0.00

Target Compounds Qvalue

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

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