

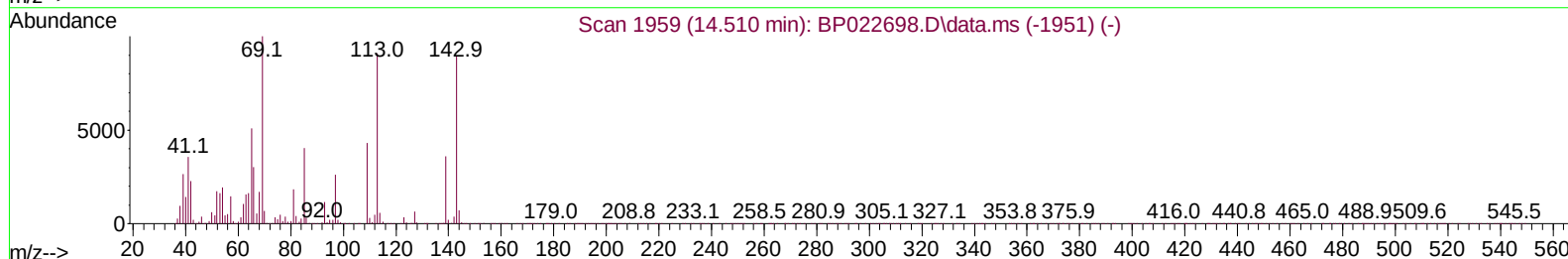
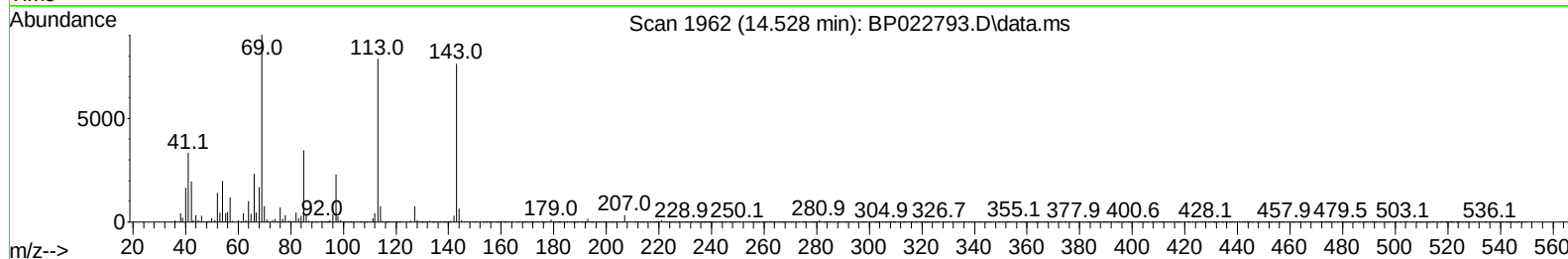
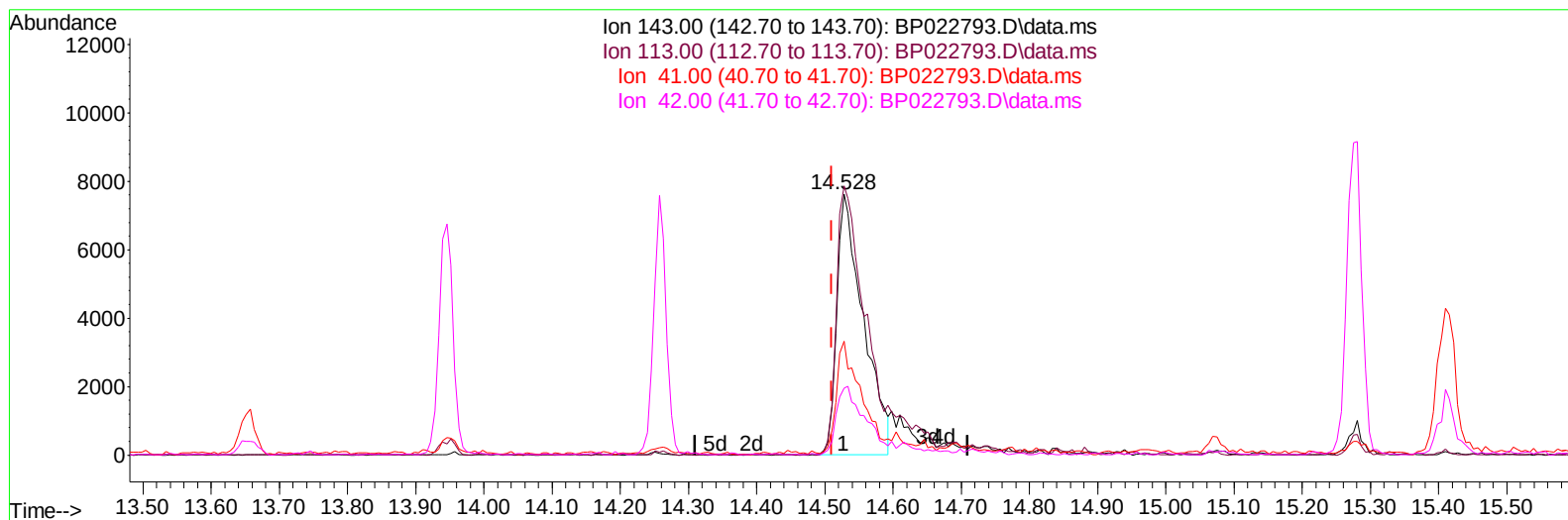
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022793.D
Acq On : 01 Nov 2024 07:15
Operator : RC/JU
Sample : P4630-13
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP07

Manual IntegrationsAPPROVED

Quant Time: Nov 01 08:02:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022793.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 4.94 ng/ul

response 20644

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.06
41.00	35.90	43.61#
42.00	23.70	25.68

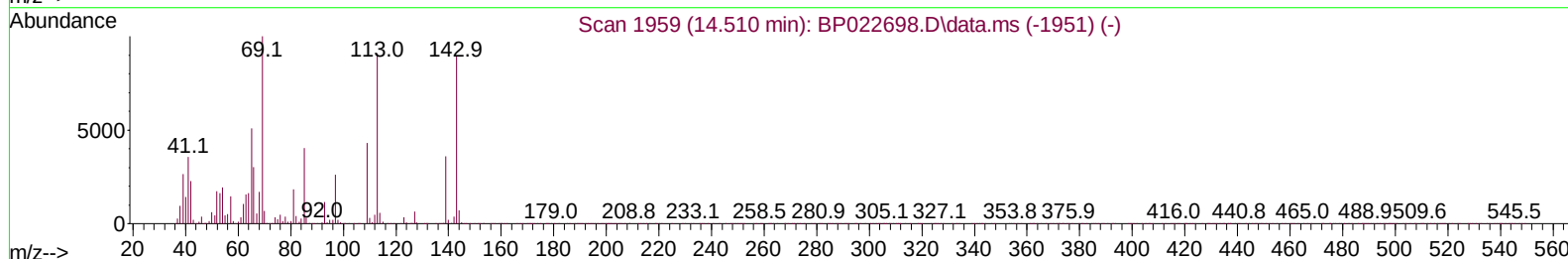
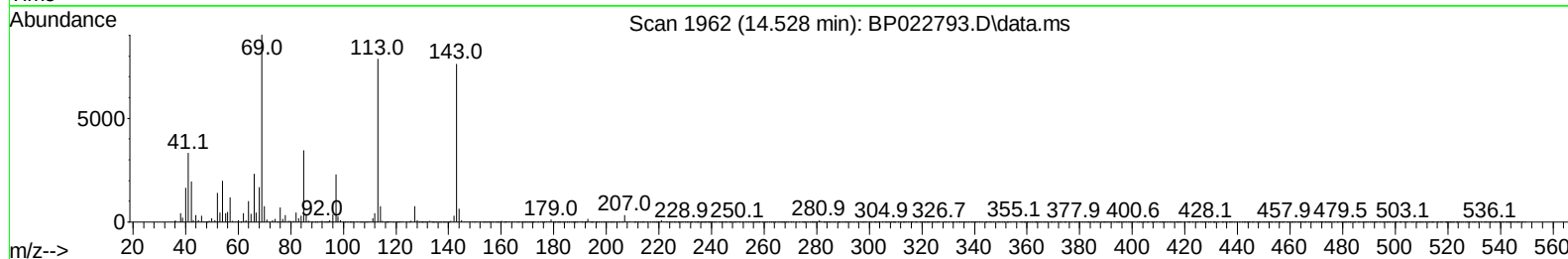
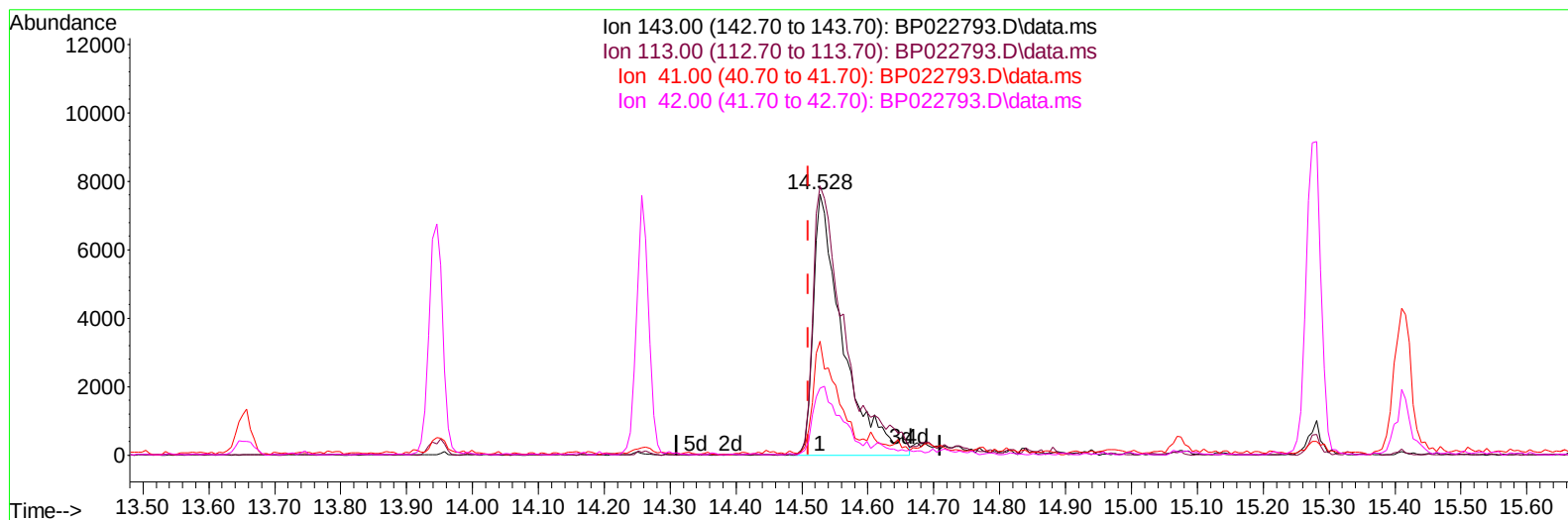
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(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 5.61 ng/ul m

response 23478

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.06
41.00	35.90	43.61#
42.00	23.70	25.68

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Quant Time: Nov 01 08:03:35 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	104491	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	392496	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	313785	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	748922	20.000	ng/ul	0.01
79) Chrysene-d12	21.504	240	866417	20.000	ng/ul	# 0.00
88) Perylene-d12	24.757	264	1077892	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	9729	3.618	ng/uL	0.00
4) Pyridine-d5	3.599	84	23958	3.246	ng/ul	0.00
7) Phenol-d5	6.834	99	85953	9.772	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	130779	25.297	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	155562	24.925	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	143804	20.487	ng/ul	-0.01
21) Nitrobenzene-d5	8.775	128	85084	28.193	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	104934	30.033	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.022	165	202500	27.269	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.528	131	146118	16.652	ng/ul	-0.01
46) Dimethylphthalate-d6	13.657	166	679084	27.239	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	738831	27.902	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	23478m	5.613	ng/ul	0.02
60) Fluorene-d10	15.281	176	626840	28.988	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	123409	25.055	ng/ul	0.00
73) Anthracene-d10	17.163	188	1173407	33.306	ng/ul	0.00
81) Pyrene-d10	19.551	212	1522320	33.226	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1900262	33.011	ng/ul	0.01

Target Compounds Qvalue

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

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