

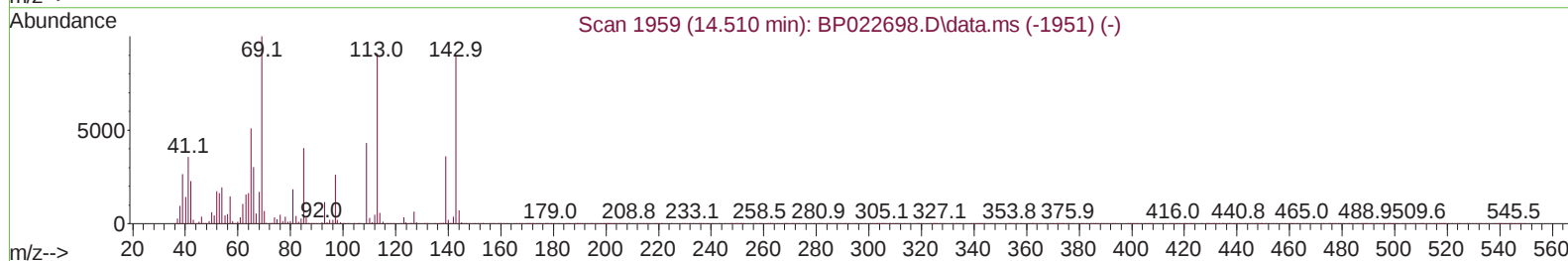
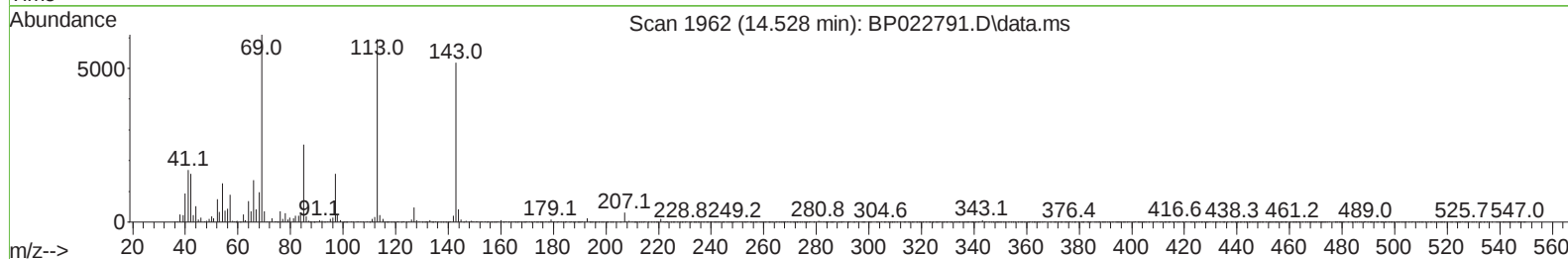
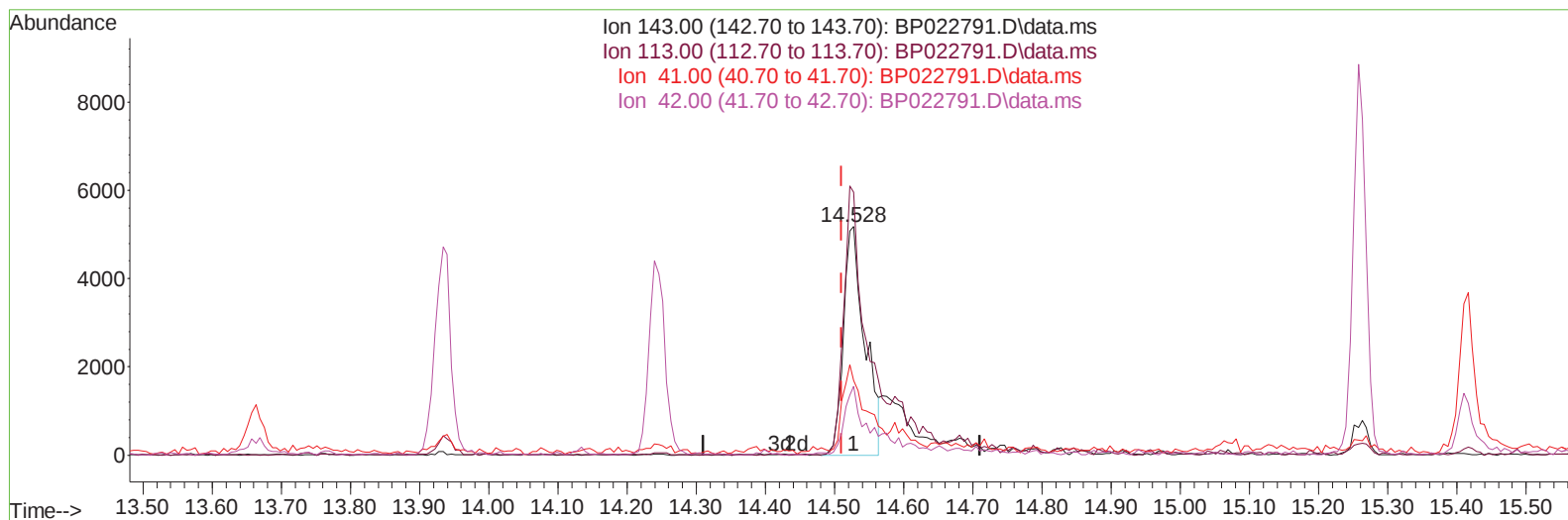
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022791.D
Acq On : 01 Nov 2024 05:54
Operator : RC/JU
Sample : PB164516BL
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK516

Manual IntegrationsAPPROVED

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

Quant Time: Nov 01 07:58:45 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



TIC: BP022791.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 3.62 ng/u1

response 11141

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	115.08
41.00	35.90	32.66
42.00	23.70	30.22#

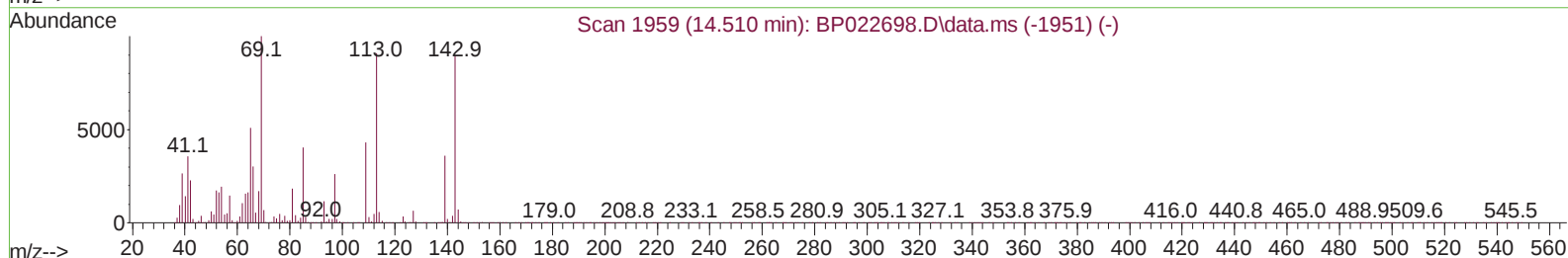
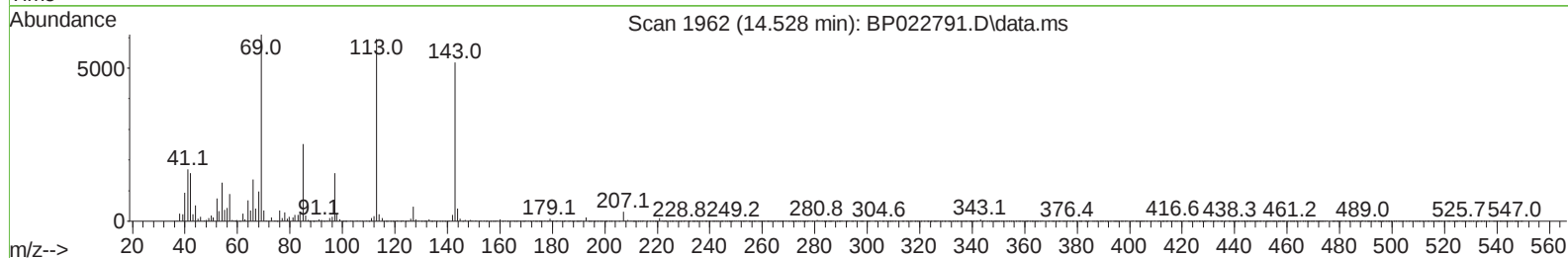
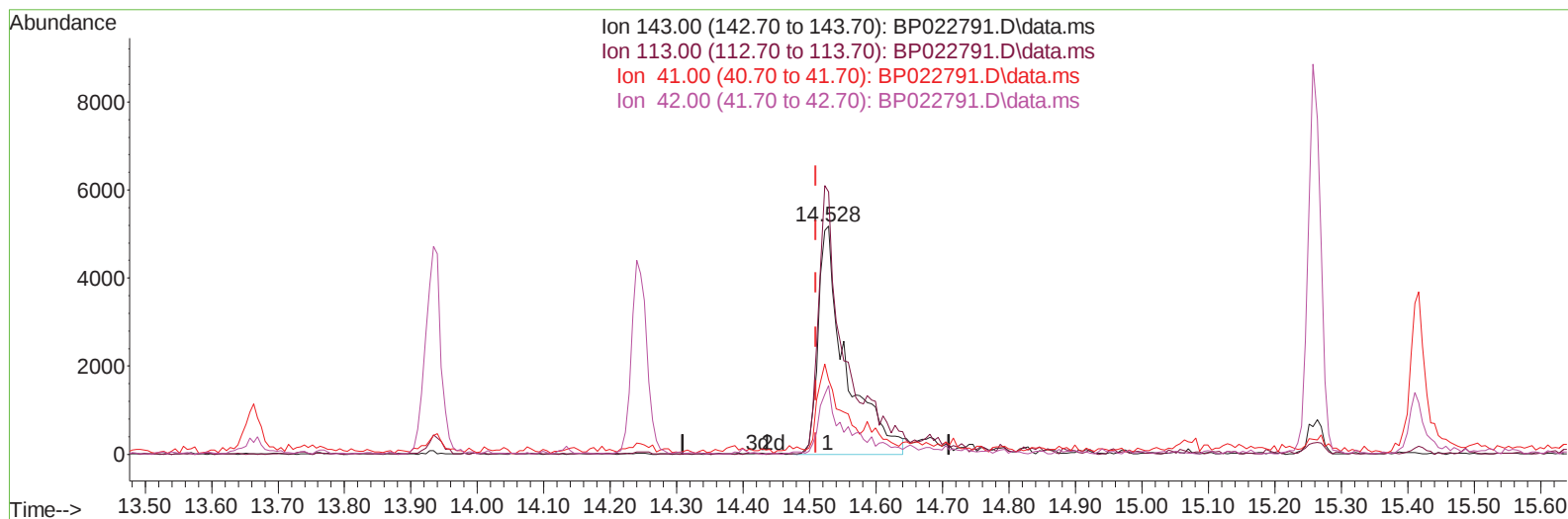
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TIC: BP022791.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 4.85 ng/ul m

response 14920

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	115.08
41.00	35.90	32.66
42.00	23.70	30.22#

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Manual IntegrationsAPPROVED

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

Quant Time: Nov 04 03:48:34 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	84038	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	312816	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.246	164	230660	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.057	188	572972	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	686222	20.000	ng/ul	0.00
88) Perylene-d12	24.769	264	847449	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	6928	3.204	ng/uL	0.00
4) Pyridine-d5	3.599	84	43815	7.380	ng/ul	0.00
7) Phenol-d5	6.828	99	46153	6.524	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	109246	26.275	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	122553	24.415	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	94113	16.671	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	71210	29.606	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	84639	30.395	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.022	165	169194	28.588	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.534	131	146325	20.923	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	558113	30.454	ng/ul	0.00
49) Acenaphthylene-d8	13.934	160	580717	29.834	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.528	143	14920m	4.853	ng/ul	0.02
60) Fluorene-d10	15.263	176	492747	30.999	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.416	200	91991	24.411	ng/ul	0.01
73) Anthracene-d10	17.169	188	895545	33.225	ng/ul	0.00
81) Pyrene-d10	19.551	212	1178099	32.465	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1509524	33.354	ng/ul	0.01

Target Compounds Qvalue

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

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