

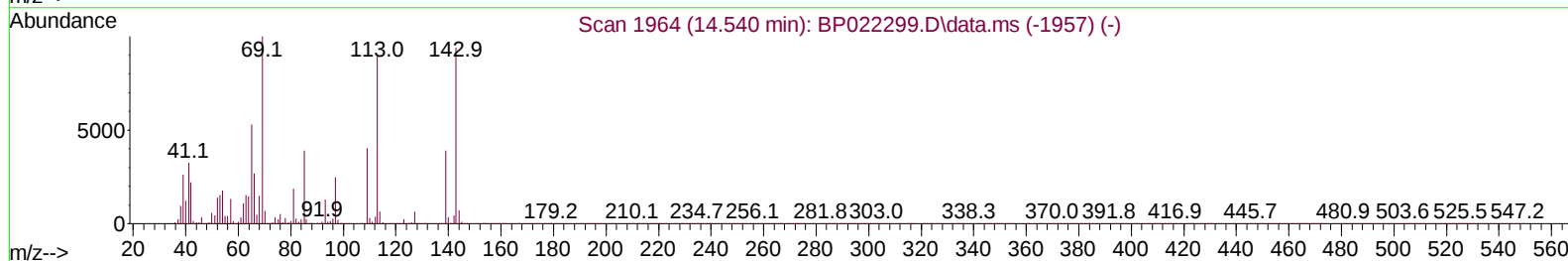
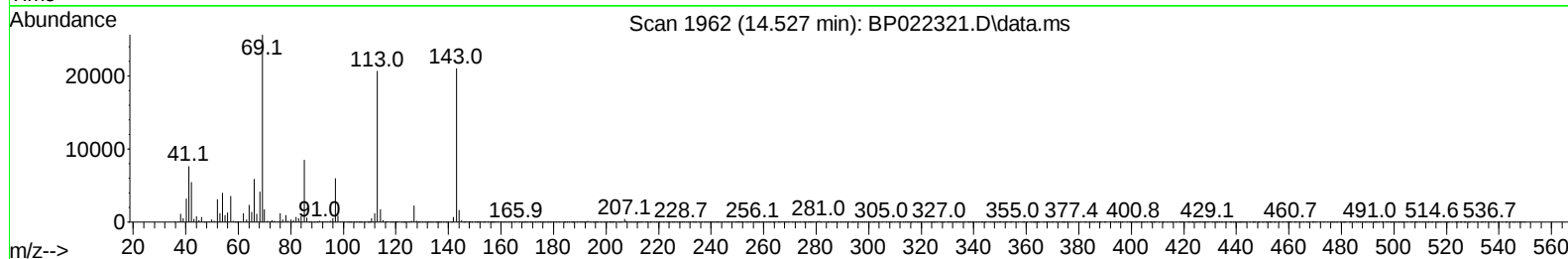
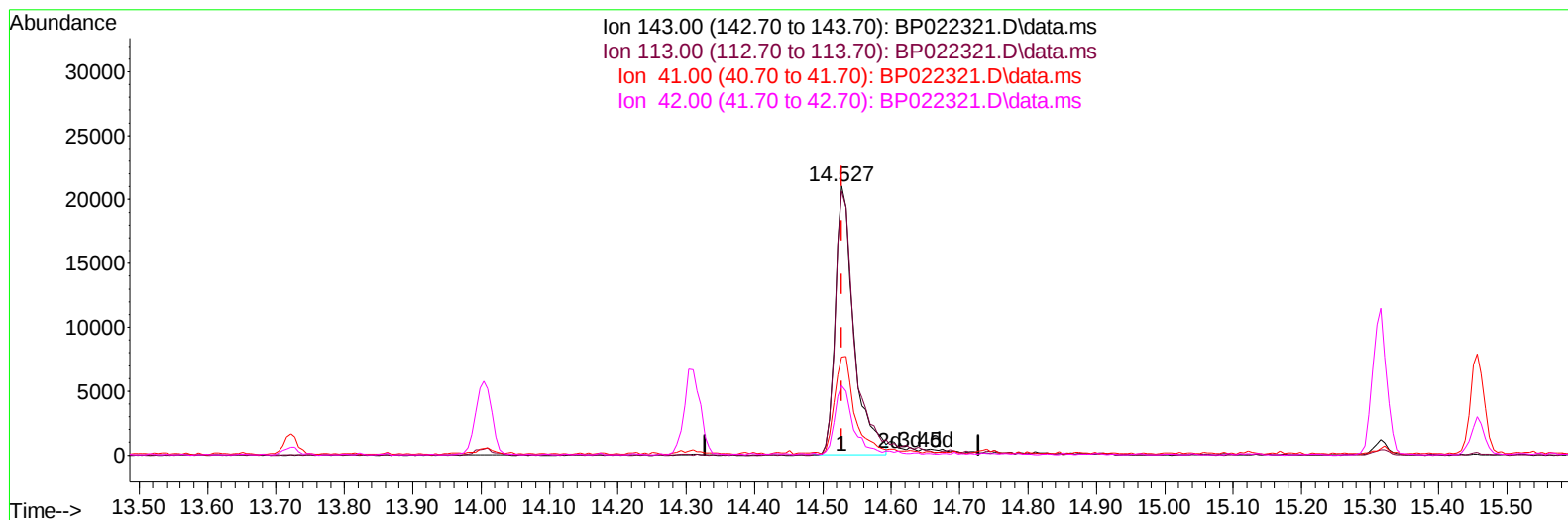
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022321.D
Acq On : 10 Oct 2024 13:29
Operator : RC/JU
Sample : P4187-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY056

Manual IntegrationsAPPROVED

Quant Time: Oct 10 14:30:39 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 08 02:14:47 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/11/2024
Supervised By :mohammad ahmed 10/11/2024



TIC: BP022321.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.527min (-0.000) 7.51 ng/u1

response 39836

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	98.21
41.00	35.90	36.54
42.00	23.70	25.88

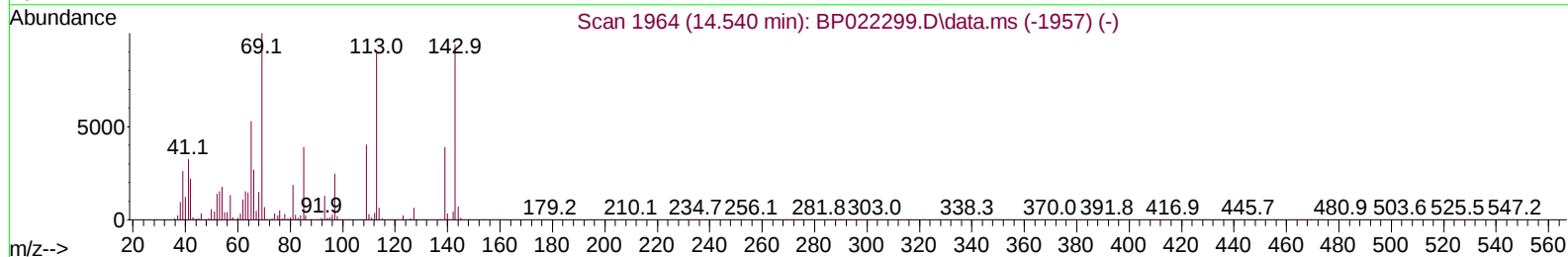
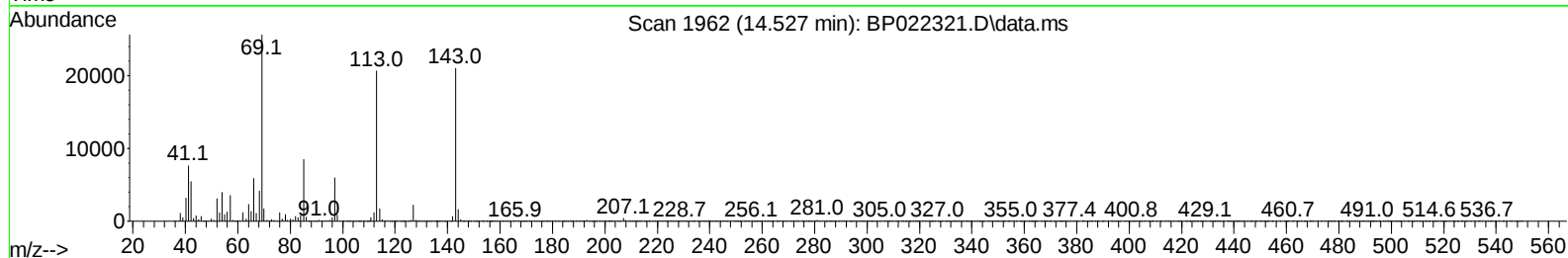
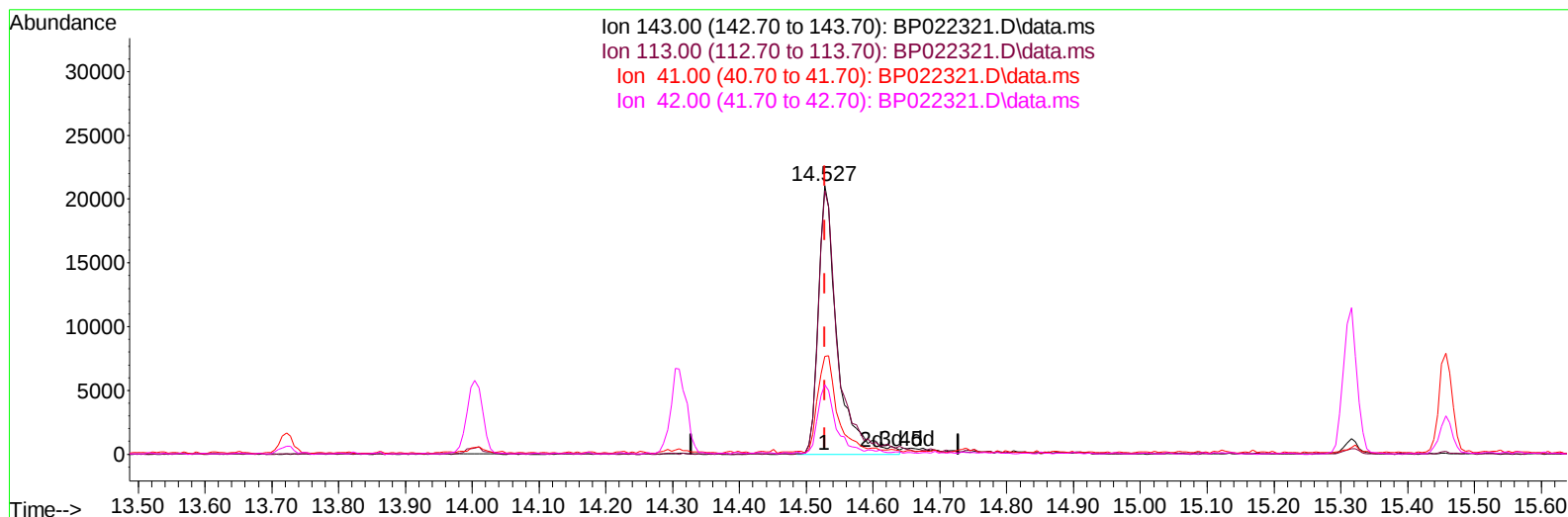
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TIC: BP022321.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.527min (-0.000) 7.86 ng/u1 m

response 41718

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	98.21
41.00	35.90	36.54
42.00	23.70	25.88

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Reviewed By :Yogesh Patel 10/11/2024
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Quant Time: Oct 11 01:46:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 02:14:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	122414	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	472619	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	398129	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	1011703	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1119018	20.000	ng/ul	0.01
88) Perylene-d12	24.839	264	1232657	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	10666	3.386	ng/uL	0.00
4) Pyridine-d5	3.646	84	61560	7.119	ng/ul	0.01
7) Phenol-d5	6.893	99	80976	7.858	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	153362	25.322	ng/ul	0.00
11) 2-Chlorophenol-d4	7.251	132	178129	24.362	ng/ul	0.00
15) 4-Methylphenol-d8	8.404	113	150413	18.291	ng/ul	0.00
21) Nitrobenzene-d5	8.845	128	100210	27.576	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	120840	28.722	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.092	165	239644	26.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.598	131	253847	24.025	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	868261	27.449	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	903212	26.884	ng/ul	0.00
54) 4-Nitrophenol-d4	14.527	143	41718m	7.861	ng/ul	0.00
60) Fluorene-d10	15.316	176	795938	29.010	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	185485	27.876	ng/ul	0.00
73) Anthracene-d10	17.216	188	1582743	33.256	ng/ul	0.00
81) Pyrene-d10	19.586	212	2033392	34.362	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.609	264	2242374	34.063	ng/ul	0.01

Target Compounds Qvalue

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

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