

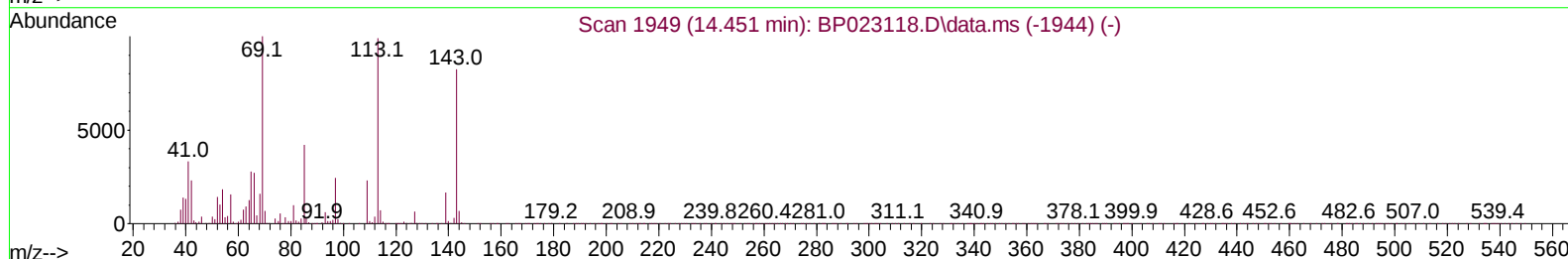
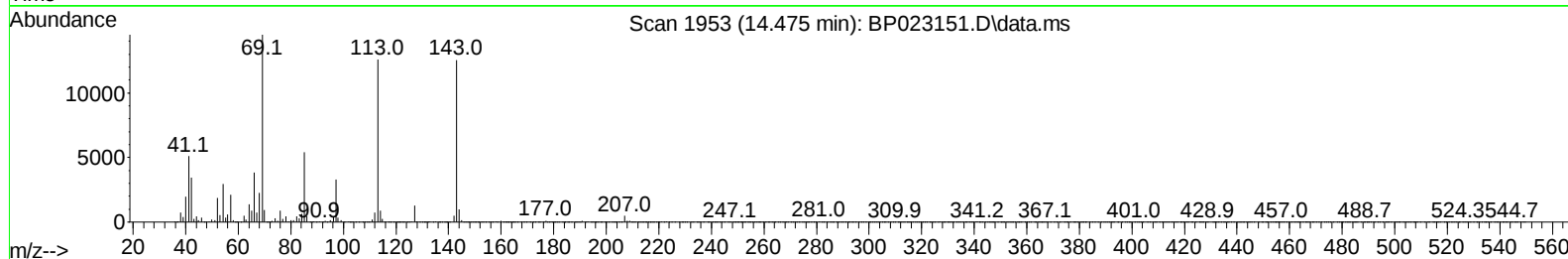
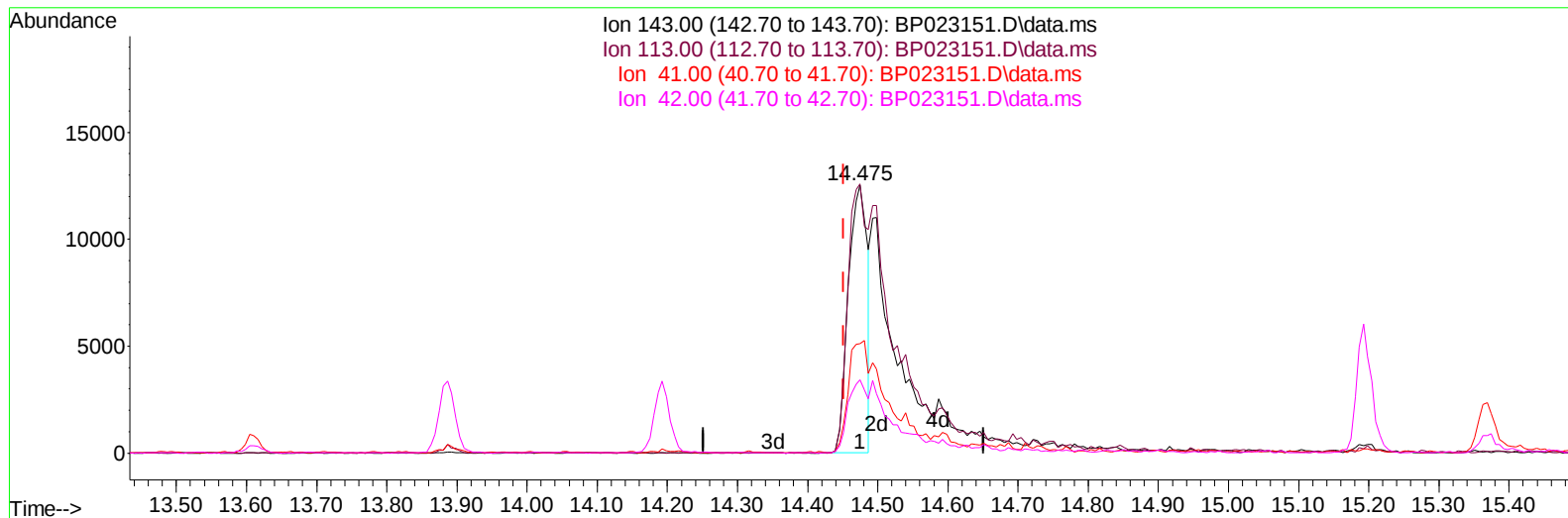
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023151.D
Acq On : 15 Nov 2024 23:34
Operator : RC/JU
Sample : PB164976BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK976

Manual IntegrationsAPPROVED

Quant Time: Nov 16 00:07:25 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023151.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.024) 13.24 ng/u1

response 23680

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	100.28
41.00	39.20	40.74
42.00	25.90	27.29

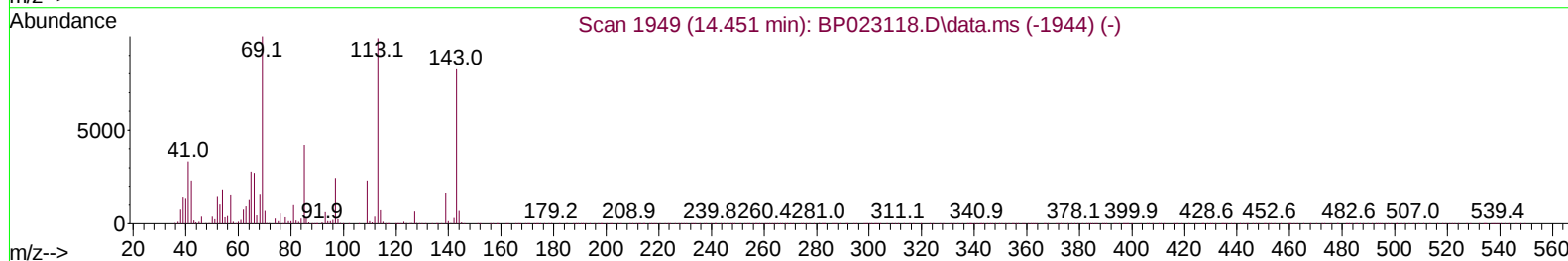
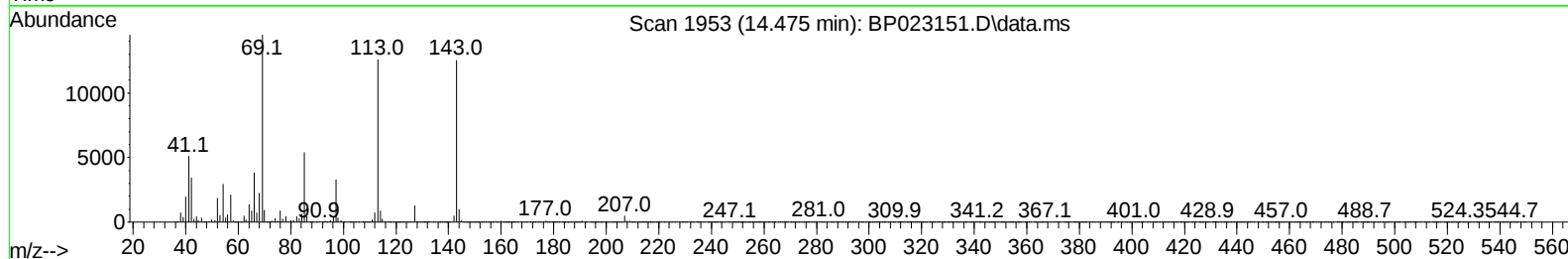
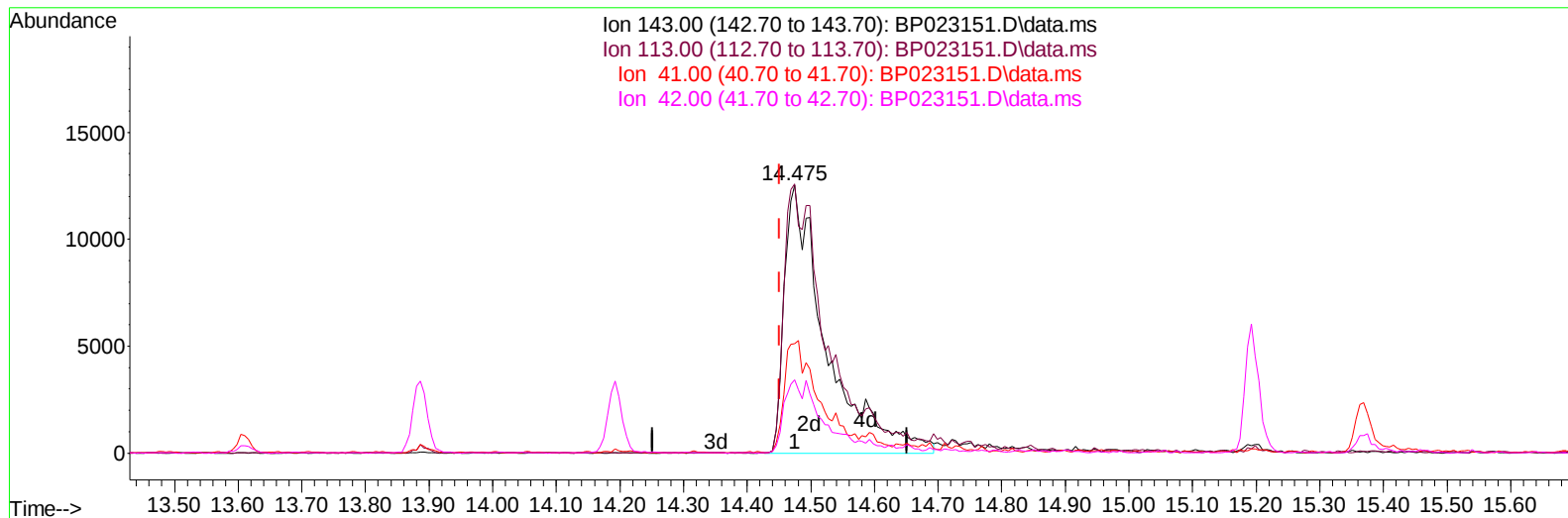
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TIC: BP023151.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.024) 31.94 ng/ul m

response 57118

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	100.28
41.00	39.20	40.74
42.00	25.90	27.29

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2024
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Quant Time: Nov 16 00:26:27 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	56639	20.000	ng/ul	0.00
20) Naphthalene-d8	10.322	136	217810	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.193	164	165120	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.992	188	420440	20.000	ng/ul	-0.02
79) Chrysene-d12	21.439	240	507682	20.000	ng/ul	0.00
88) Perylene-d12	24.663	264	554372	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	9086	7.711	ng/uL	0.00
4) Pyridine-d5	3.546	84	107616	30.740	ng/ul	0.00
7) Phenol-d5	6.775	99	137170	32.757	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	81566	33.154	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	104387	34.038	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	111592	32.384	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	53050	34.986	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	62937	34.815	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	121424	32.074	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	140963	31.211	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	434192	36.485	ng/ul	0.00
49) Acenaphthylene-d8	13.887	160	454747	36.621	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.475	143	57118m	31.937	ng/ul	0.02
60) Fluorene-d10	15.198	176	385434	36.941	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.369	200	67749	26.899	ng/ul	0.00
73) Anthracene-d10	17.104	188	671484	37.720	ng/ul	-0.01
81) Pyrene-d10	19.492	212	883854	37.797	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.427	264	1035698	39.221	ng/ul	-0.02

Target Compounds Qvalue

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

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