

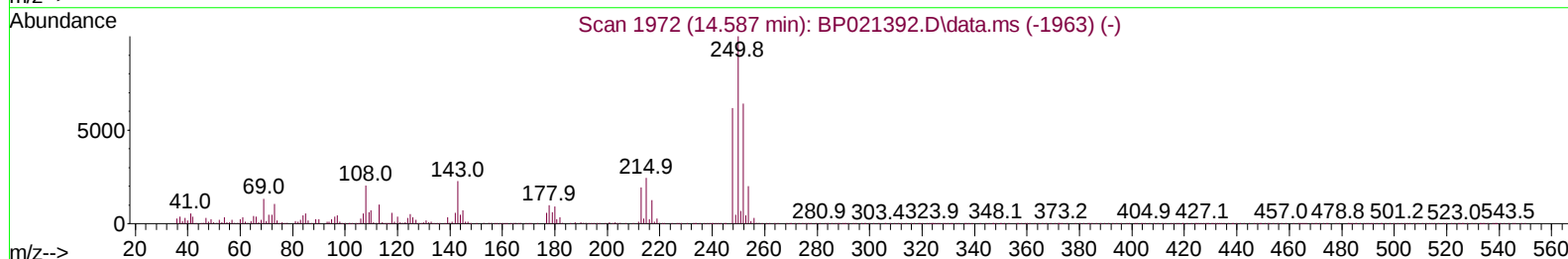
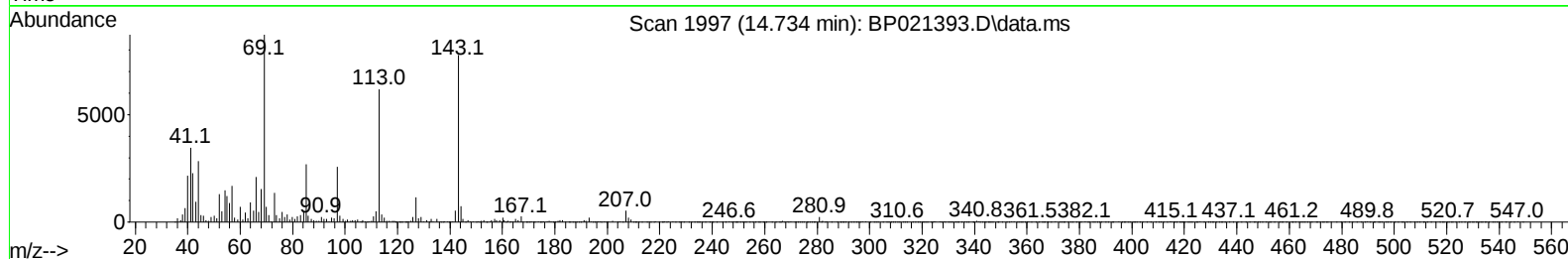
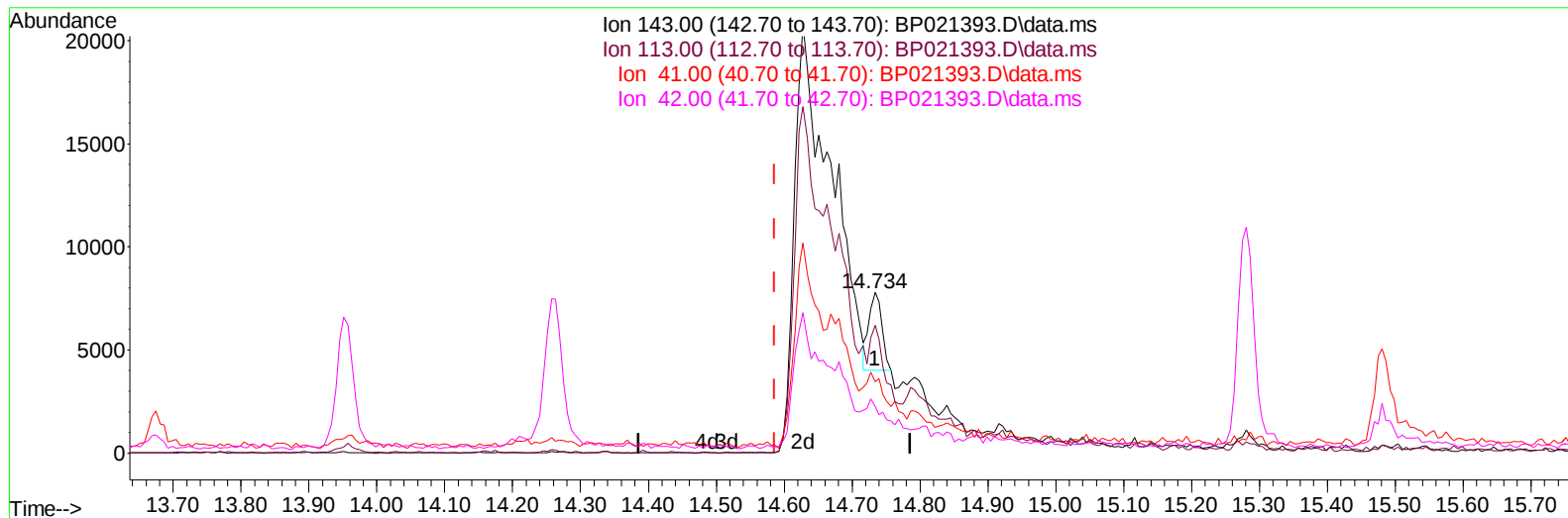
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021393.D
Acq On : 06 Aug 2024 16:11
Operator : CG/JU
Sample : PB162328BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK328

Manual IntegrationsAPPROVED

Quant Time: Aug 06 16:52:02 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/07/2024
Supervised By :mohammad ahmed 08/09/2024



TIC: BP021393.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.147) 1.19 ng/u1

response 5011

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.54
41.00	37.90	44.57
42.00	27.10	29.26

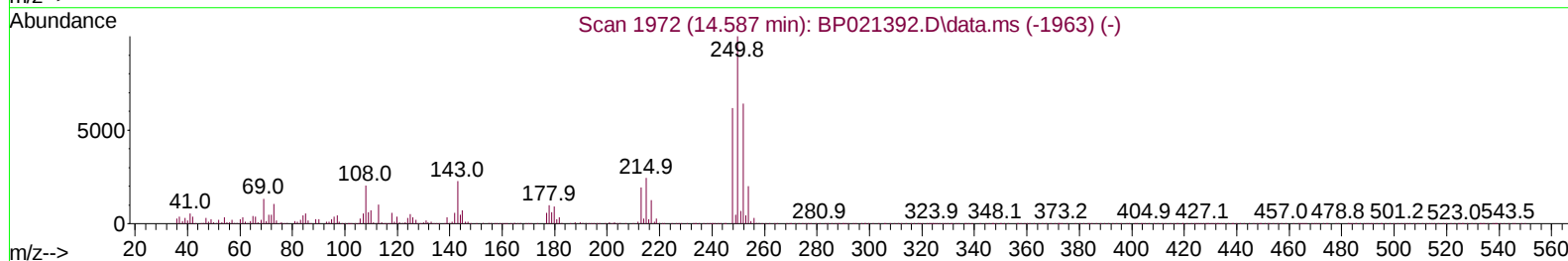
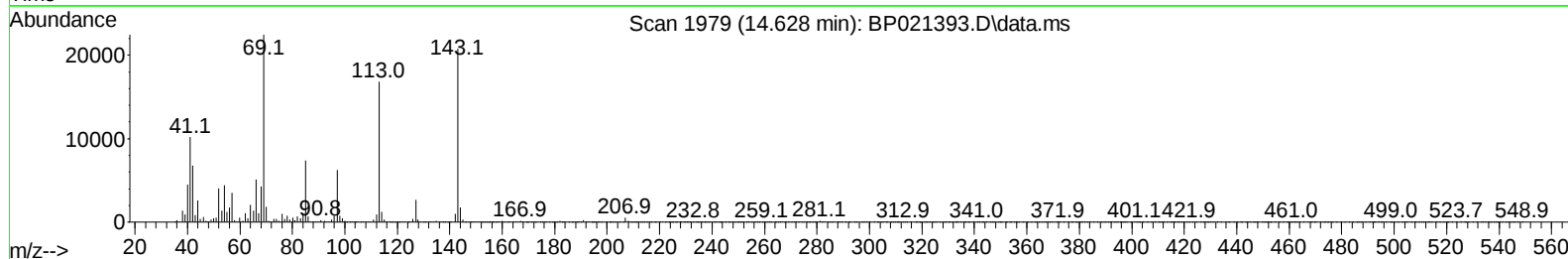
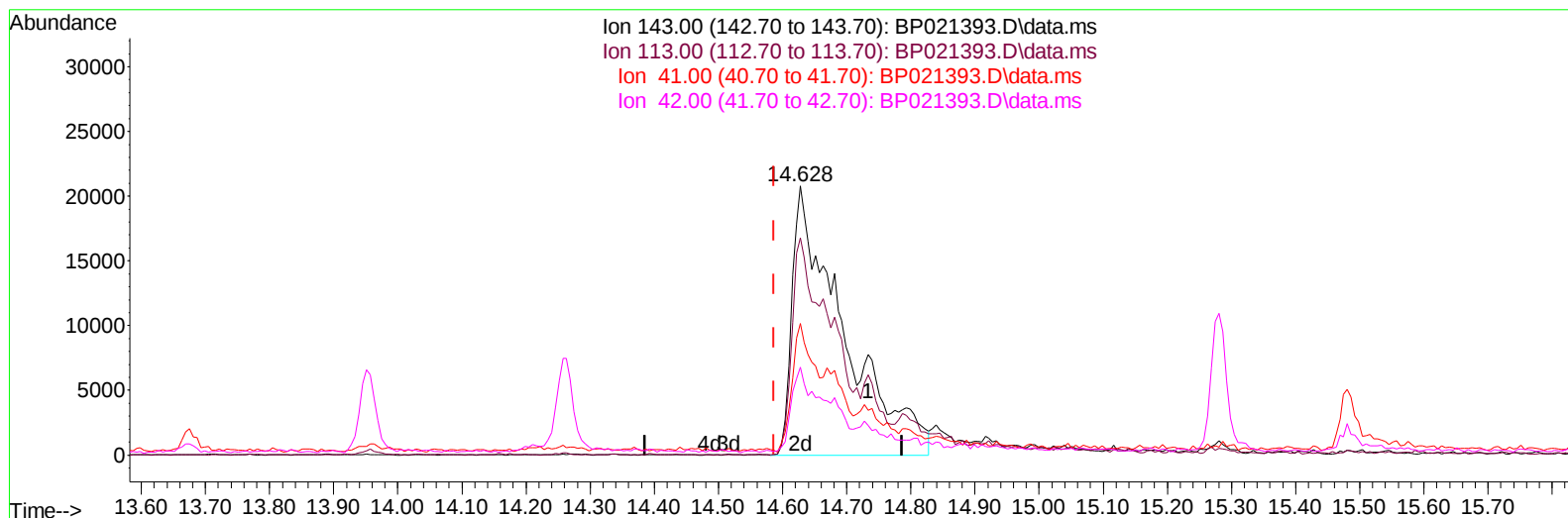
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(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.042) 27.22 ng/ul m

response 114496

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	80.86
41.00	37.90	49.03#
42.00	27.10	32.74#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/07/2024
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Quant Time: Aug 06 16:53:55 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	157039	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.381	136	624293	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.263	164	406646	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.081	188	896249	20.000	ng/ul	-0.01
79) Chrysene-d12	21.527	240	934639	20.000	ng/ul	0.00
88) Perylene-d12	24.815	264	1099367	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	19203	5.565	ng/uL	-0.02
4) Pyridine-d5	3.575	84	251570	25.180	ng/ul	-0.02
7) Phenol-d5	6.840	99	315834	24.669	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67	199462	25.707	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.169	132	267659	25.374	ng/ul	-0.01
15) 4-Methylphenol-d8	8.363	113	262264	24.664	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	125193	25.819	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	146257	26.353	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	238098	23.725	ng/ul	0.02
31) 4-Chloroaniline-d4	10.563	131	349604	26.296	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	839768	28.408	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	908404	27.367	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	114496m	27.223	ng/ul	0.04
60) Fluorene-d10	15.281	176	705415	29.087	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	111388	20.541	ng/ul	0.02
73) Anthracene-d10	17.181	188	1140186	28.643	ng/ul	-0.01
81) Pyrene-d10	19.569	212	1453734	25.158	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.580	264	1577391	29.092	ng/ul	-0.02

Target Compounds Qvalue

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

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