

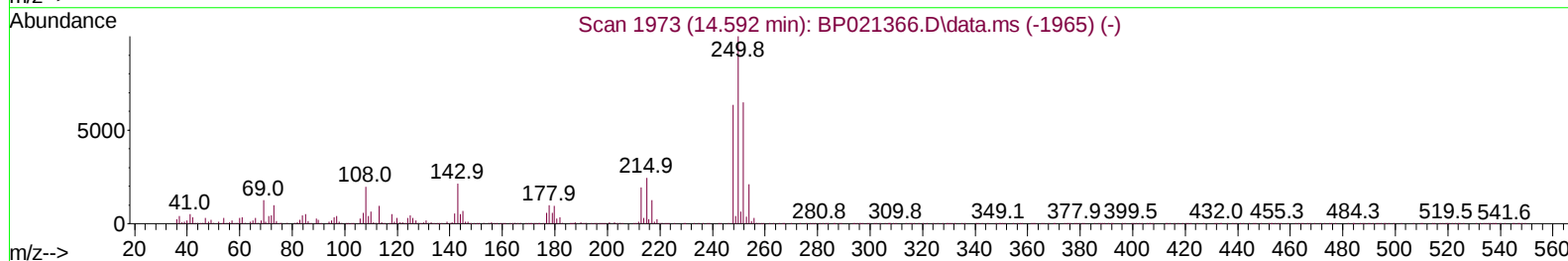
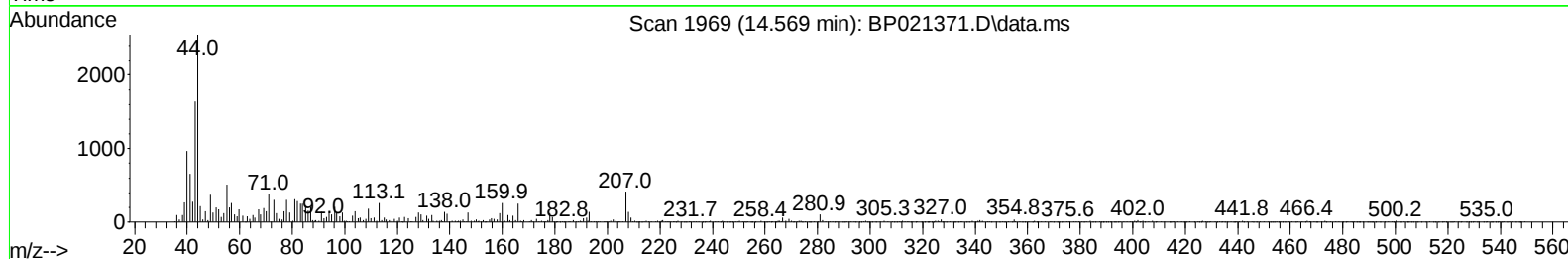
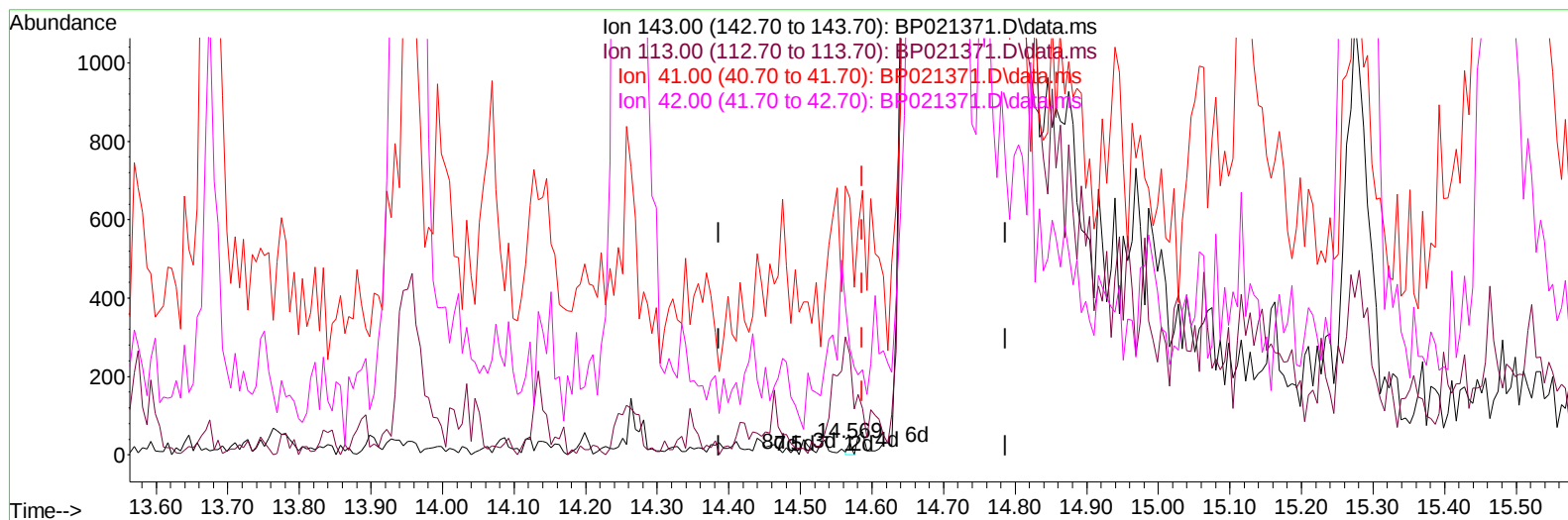
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080524\
Data File : BP021371.D
Acq On : 05 Aug 2024 13:49
Operator : CG/JU
Sample : P3300-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G91

Manual IntegrationsAPPROVED

Quant Time: Aug 05 14:32:08 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 :Jagrut Upadhyay 08/06/2024
Supervised By :mohammad ahmed 08/07/2024



TIC: BP021371.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.017) 0.00 ng/u1

response 11

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	800.00#
41.00	37.90	2046.88#
42.00	27.10	853.13#

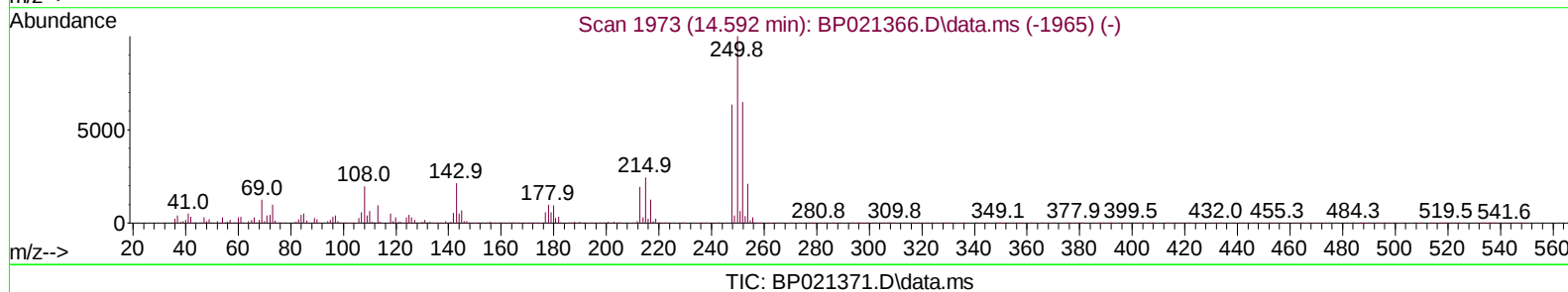
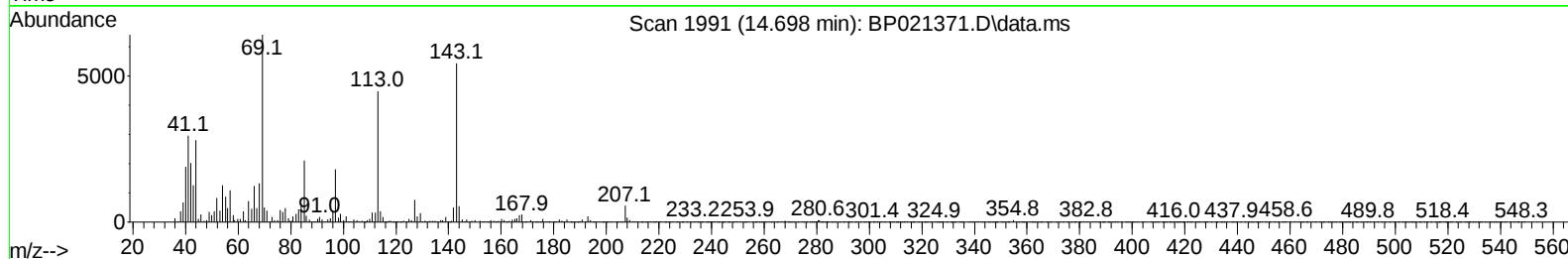
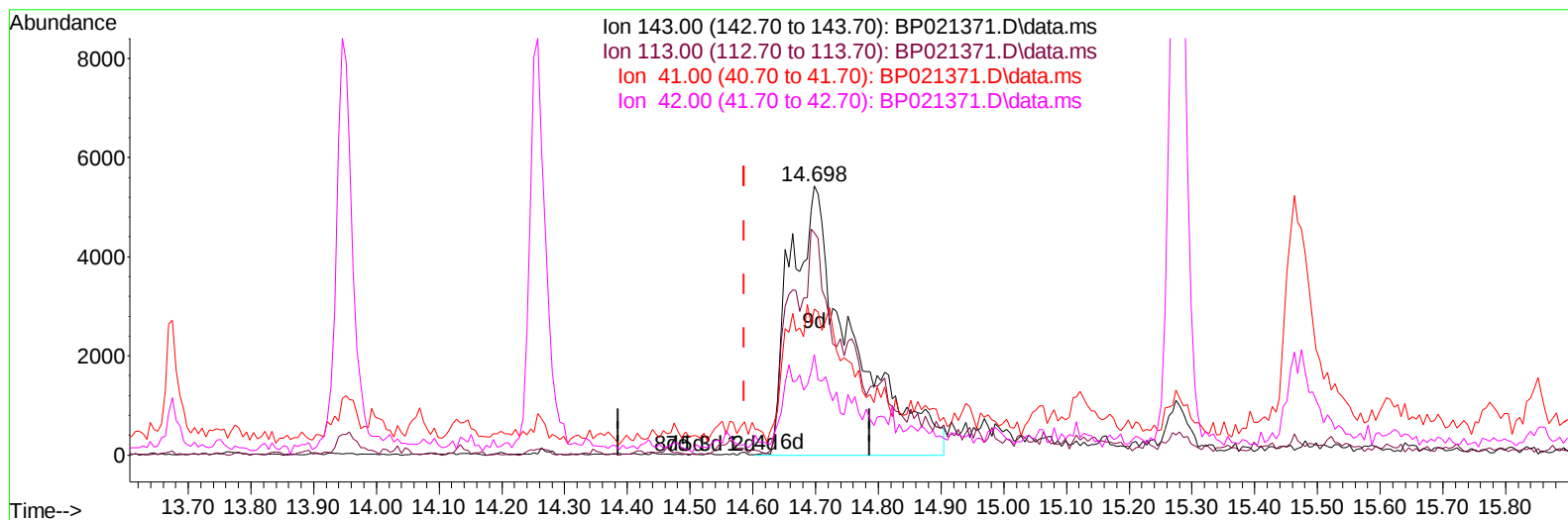
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(54) 4-Nitrophenol-d4 (S)

14.698min (+ 0.112) 7.54 ng/ul m

response 36833

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	82.57
41.00	37.90	54.31#
42.00	27.10	37.32#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	166752	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.375	136	693611	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.257	164	472004	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.081	188	1071970	20.000	ng/ul	-0.01
79) Chrysene-d12	21.527	240	1090312	20.000	ng/ul	0.00
88) Perylene-d12	24.816	264	1280070	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	16713	4.561	ng/uL	-0.02
4) Pyridine-d5	3.587	84	61927	5.837	ng/ul	0.00
7) Phenol-d5	6.858	99	96075	7.067	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.969	67	217002	26.339	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	264403	23.605	ng/ul	-0.01
15) 4-Methylphenol-d8	8.358	113	205548	18.204	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	136326	25.305	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	155574	25.230	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.052	165	271878	24.384	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	313032	21.192	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	1007959	29.376	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1084629	28.152	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.698	143	36833m	7.545	ng/ul	0.11
60) Fluorene-d10	15.275	176	871761	30.969	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	146958	22.658	ng/ul	0.00
73) Anthracene-d10	17.186	188	1511254	31.741	ng/ul	0.00
81) Pyrene-d10	19.575	212	1877772	27.857	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	2066488	32.732	ng/ul	-0.01

Target Compounds Qvalue

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

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