

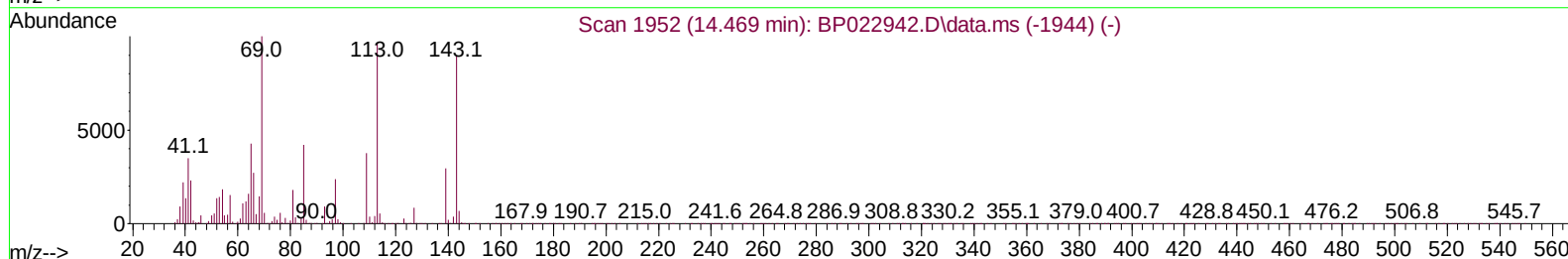
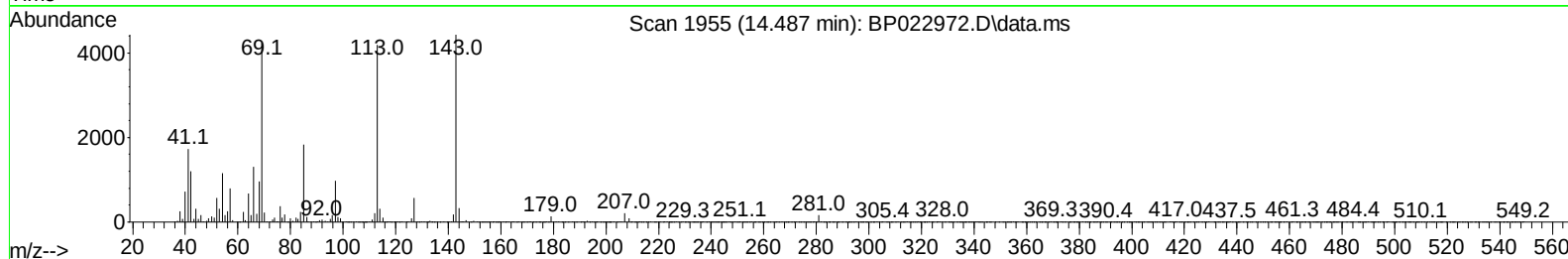
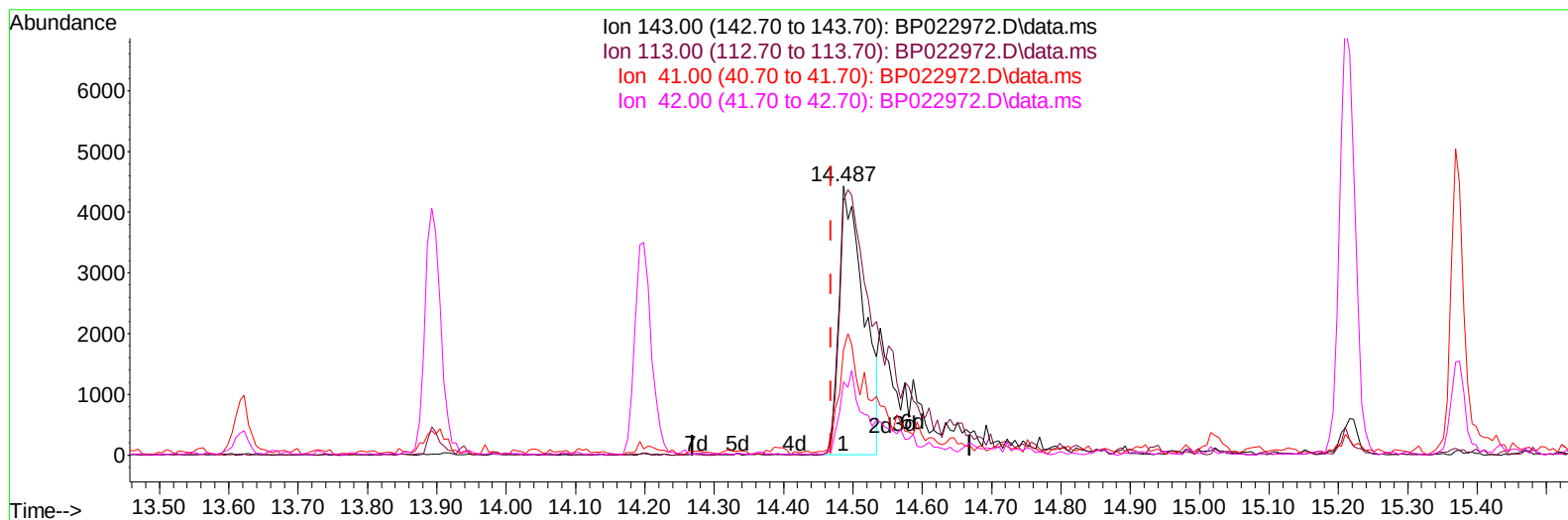
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022972.D
Acq On : 09 Nov 2024 18:36
Operator : RC/JU
Sample : P4746-03
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP78

Manual IntegrationsAPPROVED

Quant Time: Nov 09 23:24:20 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 08 23:49:48 2024
Response via : Initial Calibration

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



TIC: BP022972.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.018) 5.58 ng/ul

response 10903

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	94.74
41.00	39.20	39.11
42.00	25.90	27.16

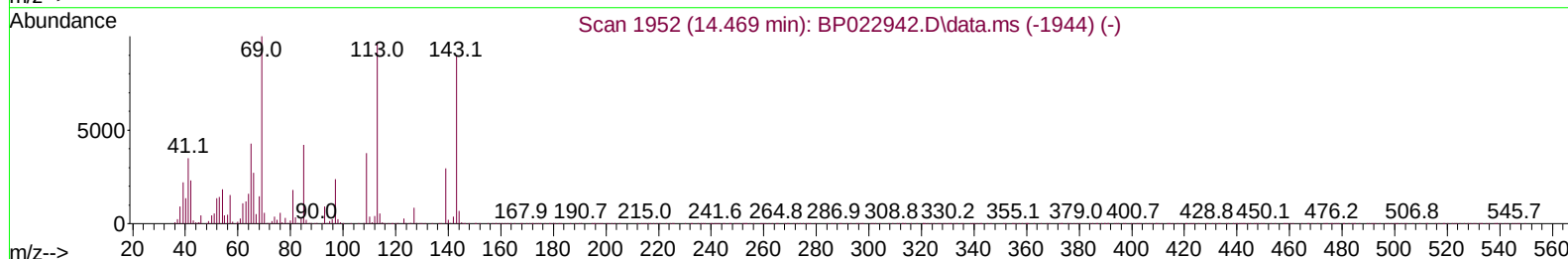
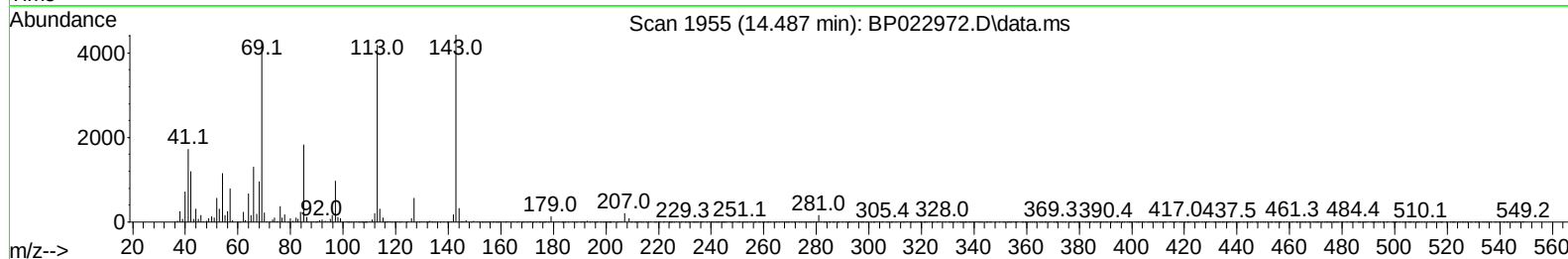
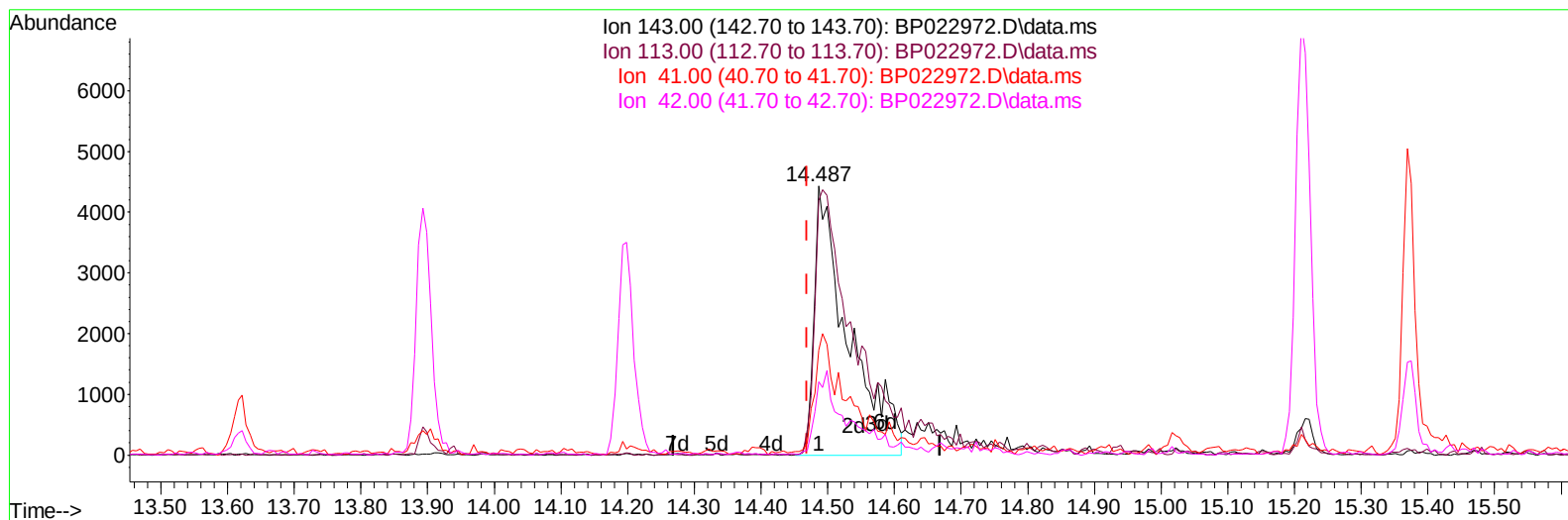
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TIC: BP022972.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.018) 8.08 ng/ul m

response 15790

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	94.74
41.00	39.20	39.11
42.00	25.90	27.16

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	56963	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	222127	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.198	164	180447	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.004	188	473724	20.000	ng/ul	-0.03
79) Chrysene-d12	21.445	240	566951	20.000	ng/ul	-0.01
88) Perylene-d12	24.668	264	681885	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	5726	4.832	ng/uL	0.00
4) Pyridine-d5	3.570	84	13031	3.701	ng/ul	0.01
7) Phenol-d5	6.799	99	37481	8.900	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	86382	34.912	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	97885	31.736	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	77391	22.331	ng/ul	-0.01
21) Nitrobenzene-d5	8.728	128	55847	36.114	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	69450	37.671	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	135336	35.054	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	97919	21.259	ng/ul	0.00
46) Dimethylphthalate-d6	13.622	166	522433	40.171	ng/ul	0.00
49) Acenaphthylene-d8	13.893	160	505422	37.245	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.487	143	15790m	8.079	ng/ul	0.02
60) Fluorene-d10	15.216	176	442241	38.785	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.375	200	99334	35.003	ng/ul	0.00
73) Anthracene-d10	17.110	188	810531	40.409	ng/ul	-0.01
81) Pyrene-d10	19.498	212	1060458	40.609	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.445	264	1389903	42.792	ng/ul	-0.02

Target Compounds Qvalue

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

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