

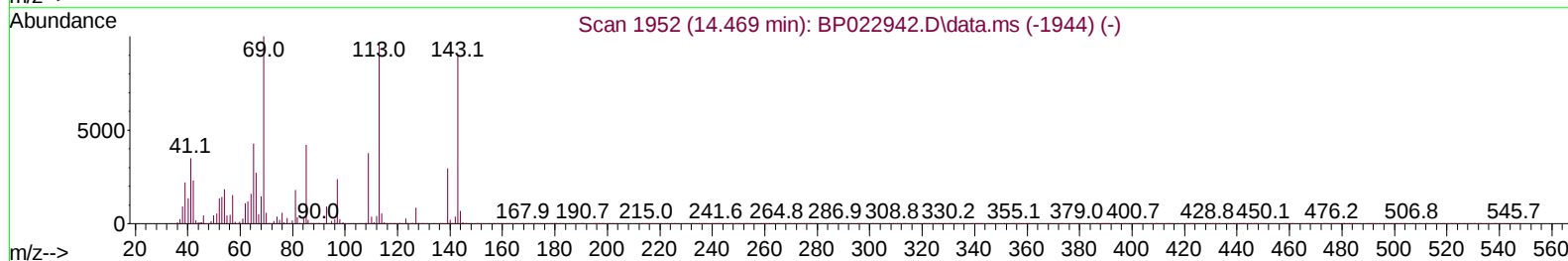
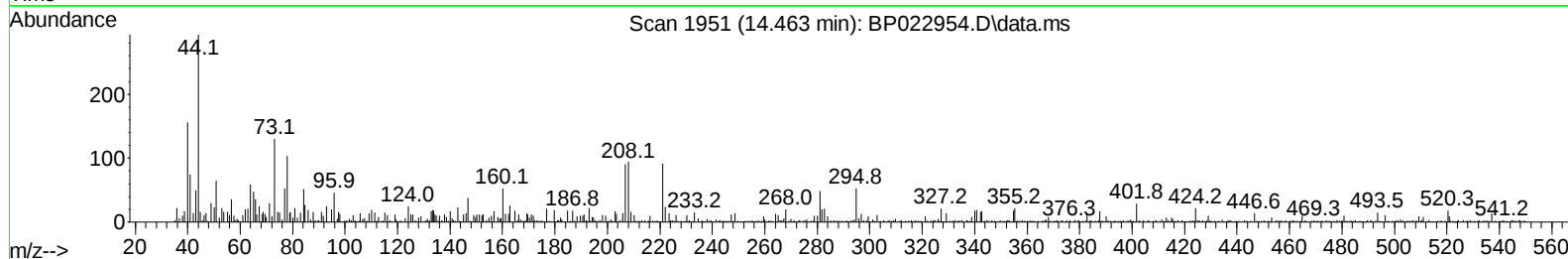
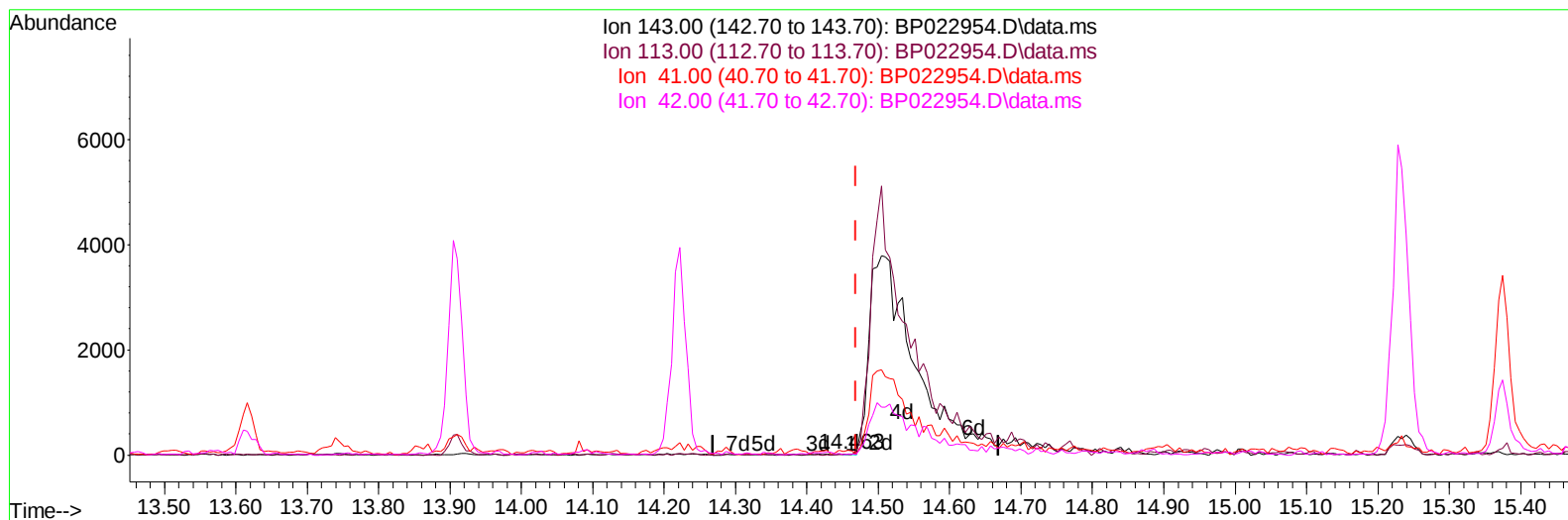
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022954.D
 Acq On : 09 Nov 2024 06:28
 Operator : RC/JU
 Sample : P4744-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP66

Manual IntegrationsAPPROVED

Quant Time: Nov 09 22:10:47 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: BP022954.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.463min (-0.006) 0.01 ng/u1

response 12

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	32.00#
41.00	39.20	300.00#
42.00	25.90	56.00#

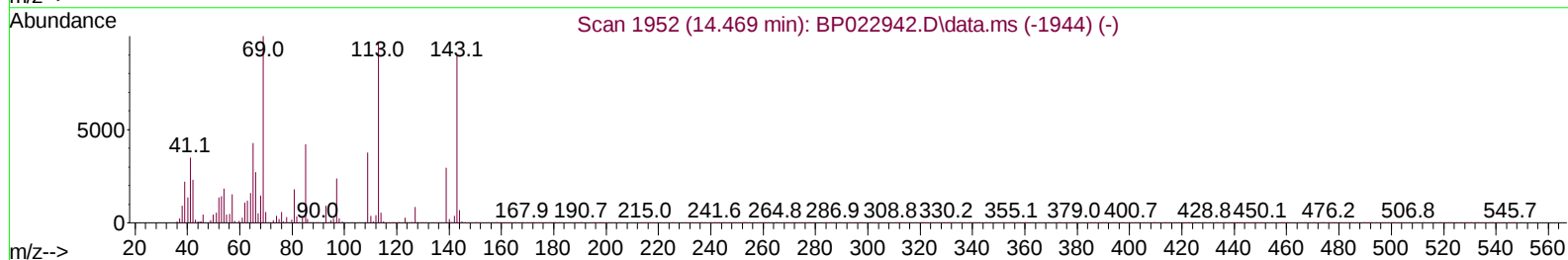
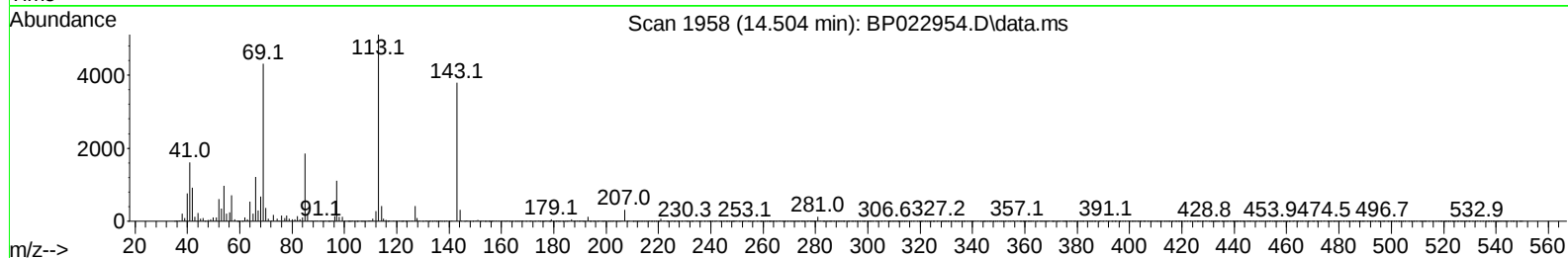
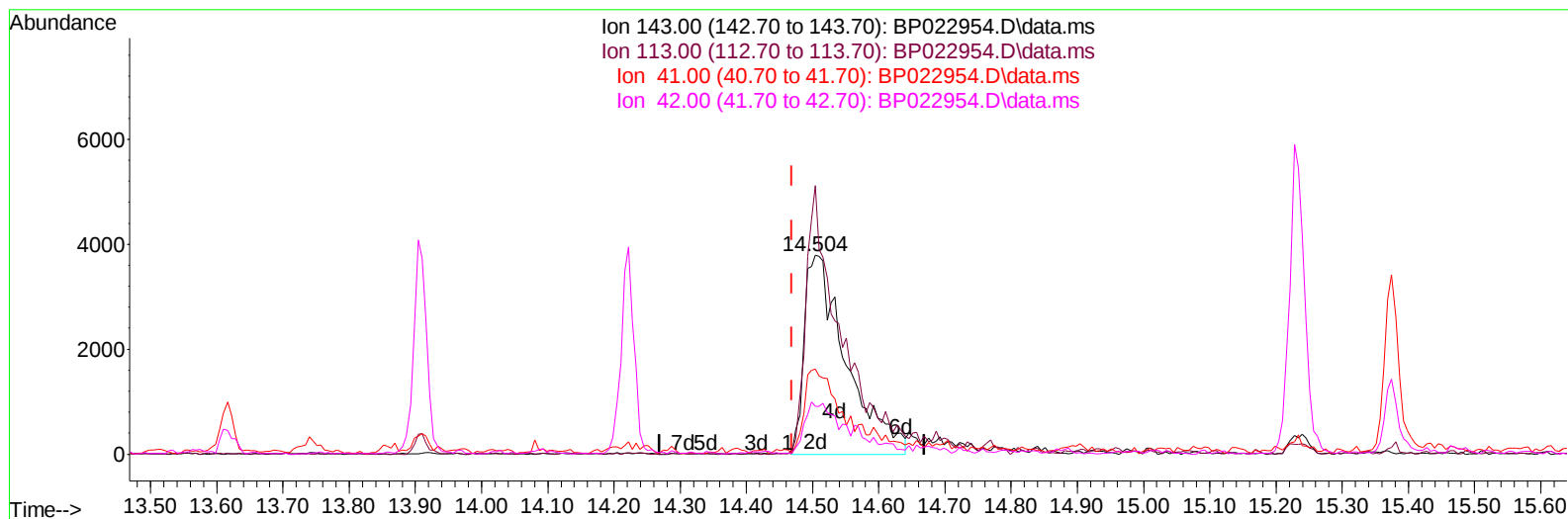
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(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.035) 9.56 ng/ul m

response 16728

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	134.91#
41.00	39.20	42.69
42.00	25.90	24.17

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	52256	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	197662	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.222	164	161496	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	404298	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	498887	20.000	ng/ul	-0.01
88) Perylene-d12	24.668	264	643136	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	5669	5.215	ng/uL	0.00
4) Pyridine-d5	3.569	84	20131	6.233	ng/ul	0.01
7) Phenol-d5	6.799	99	37486	9.703	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	79615	35.075	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	90163	31.866	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	73569	23.140	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	50576	36.754	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	59808	36.457	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	118841	34.592	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	111318	27.160	ng/ul	0.00
46) Dimethylphthalate-d6	13.616	166	458676	39.407	ng/ul	0.00
49) Acenaphthylene-d8	13.910	160	435224	35.836	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	16728m	9.563	ng/ul	0.04
60) Fluorene-d10	15.234	176	383436	37.574	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	80739	33.336	ng/ul	0.00
73) Anthracene-d10	17.122	188	659196	38.508	ng/ul	0.00
81) Pyrene-d10	19.492	212	891552	38.798	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.445	264	1231613	40.203	ng/ul	-0.02

Target Compounds Qvalue

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

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