

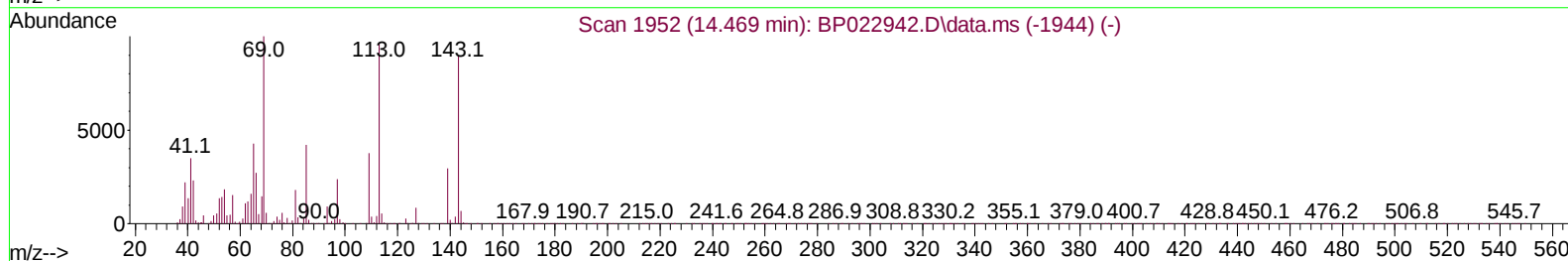
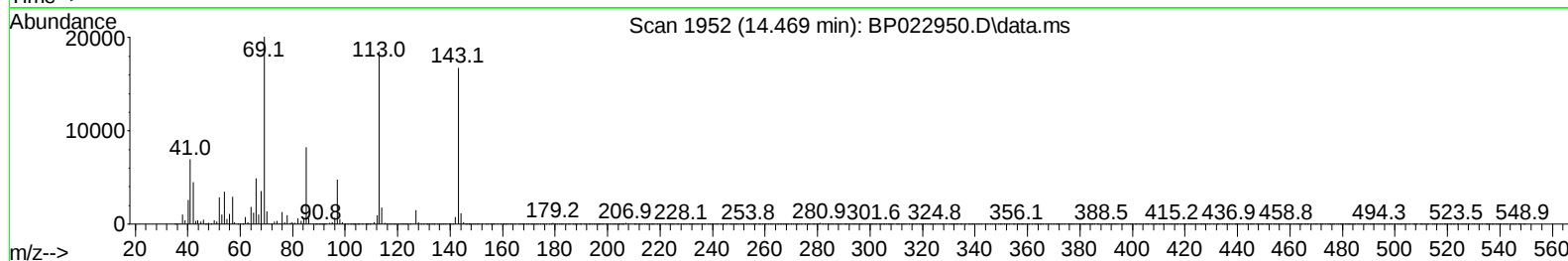
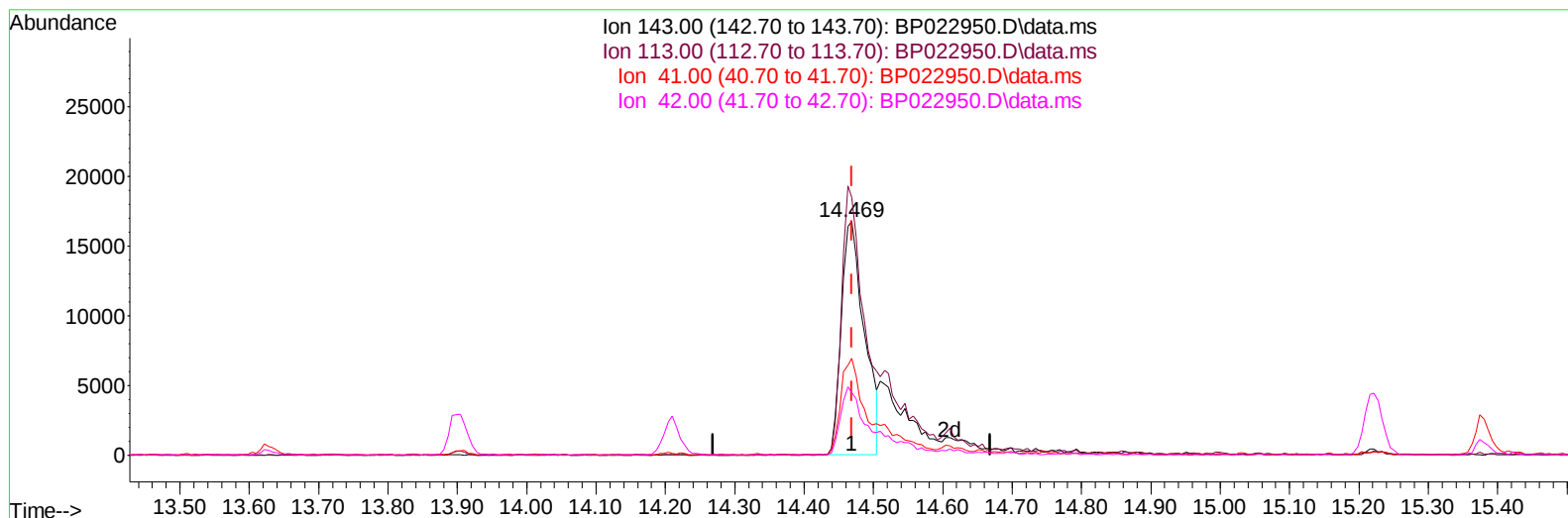
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022950.D
 Acq On : 09 Nov 2024 03:45
 Operator : RC/JU
 Sample : PB164778BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK778

Manual IntegrationsAPPROVED

Quant Time: Nov 09 04:22:37 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: BP022950.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (-0.000) 23.95 ng/u1

response 38251

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	110.30
41.00	39.20	41.36
42.00	25.90	26.66

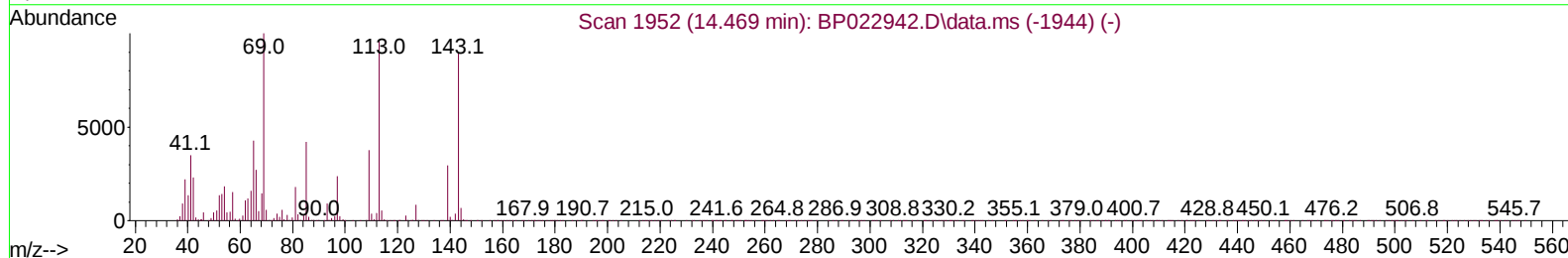
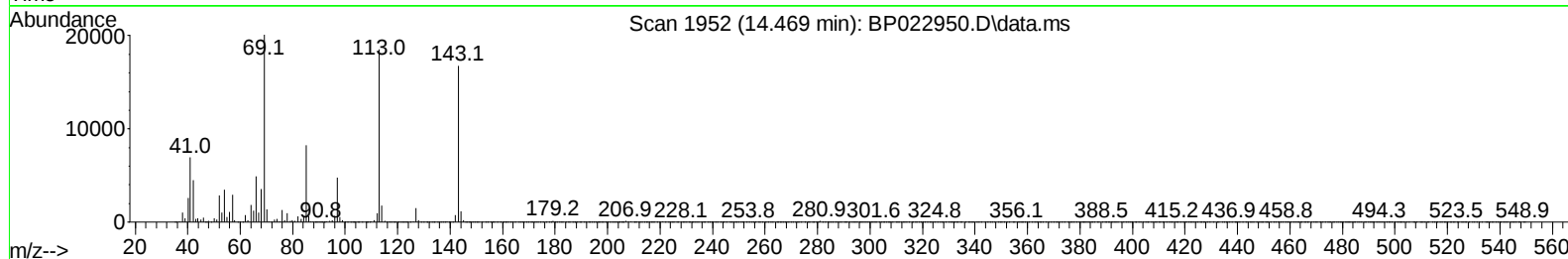
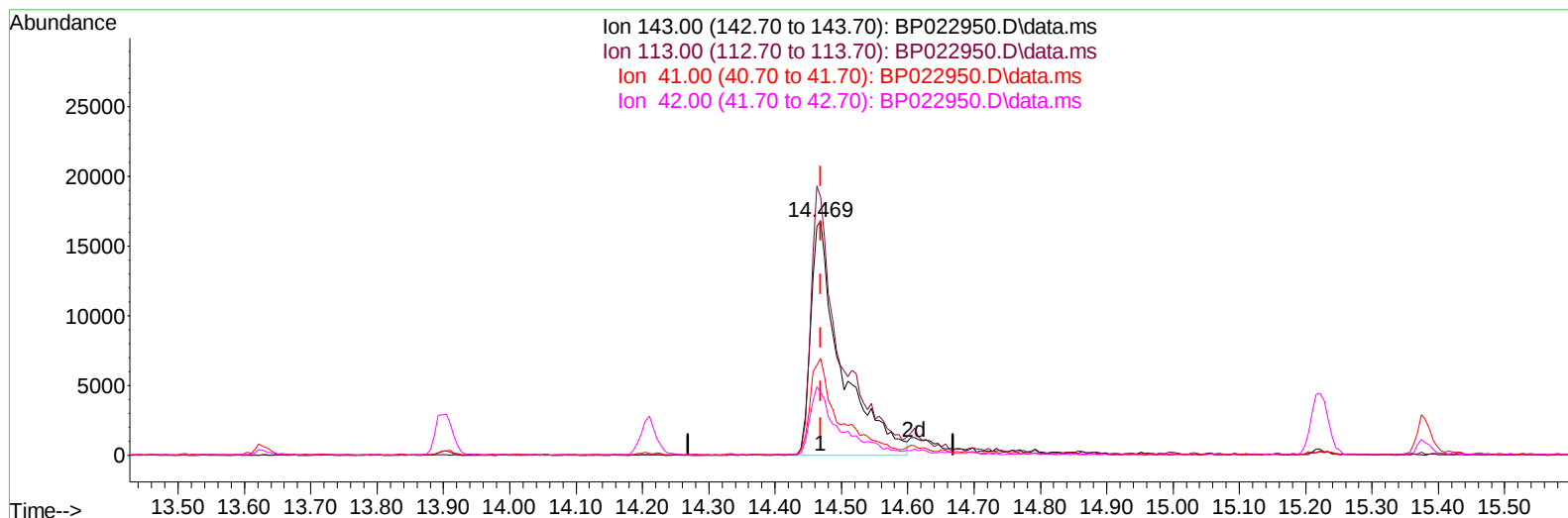
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TIC: BP022950.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (-0.000) 33.53 ng/ul m

response 53549

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	110.30
41.00	39.20	41.36
42.00	25.90	26.66

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 ALS Vial : 35 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	48648	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	183879	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.210	164	147449	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.022	188	360193	20.000	ng/ul	-0.01
79) Chrysene-d12	21.451	240	427165	20.000	ng/ul	0.00
88) Perylene-d12	24.674	264	499170	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	8209	8.111	ng/uL	0.00
4) Pyridine-d5	3.552	84	105708	35.155	ng/ul	0.00
7) Phenol-d5	6.793	99	127708	35.507	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	78144	36.980	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	97910	37.170	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	105601	35.679	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	50399	39.371	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	59712	39.127	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	112540	35.213	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	142045	37.255	ng/ul	0.00
46) Dimethylphthalate-d6	13.628	166	428206	40.294	ng/ul	0.01
49) Acenaphthylene-d8	13.898	160	435291	39.256	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	53549m	33.530	ng/ul	0.00
60) Fluorene-d10	15.222	176	363822	39.048	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.381	200	74257	34.414	ng/ul	0.01
73) Anthracene-d10	17.127	188	635975	41.701	ng/ul	0.00
81) Pyrene-d10	19.498	212	811538	41.246	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.451	264	1026647	43.177	ng/ul	-0.02

Target Compounds Qvalue

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

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