

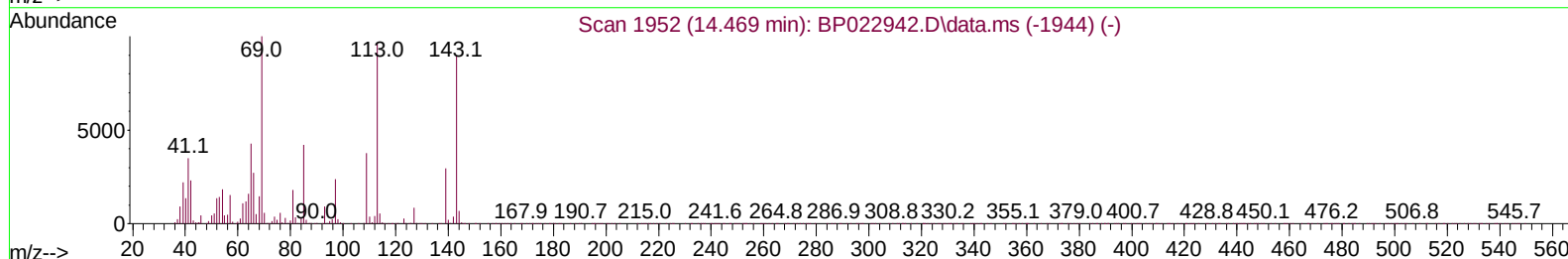
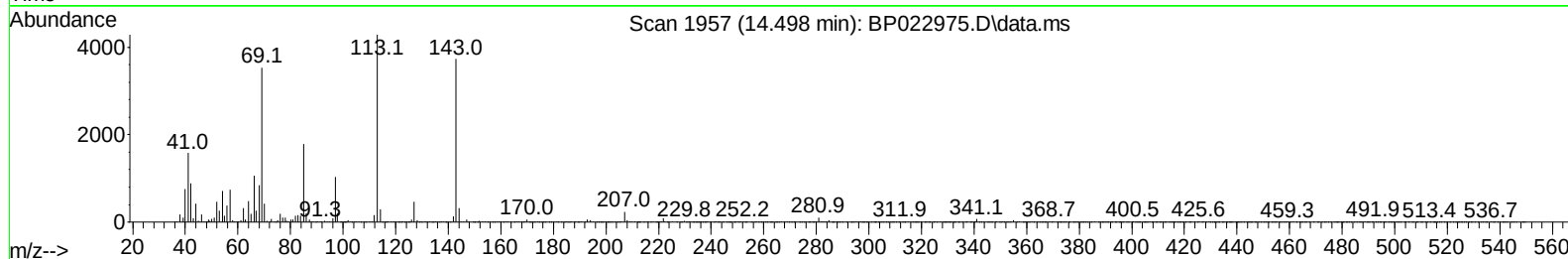
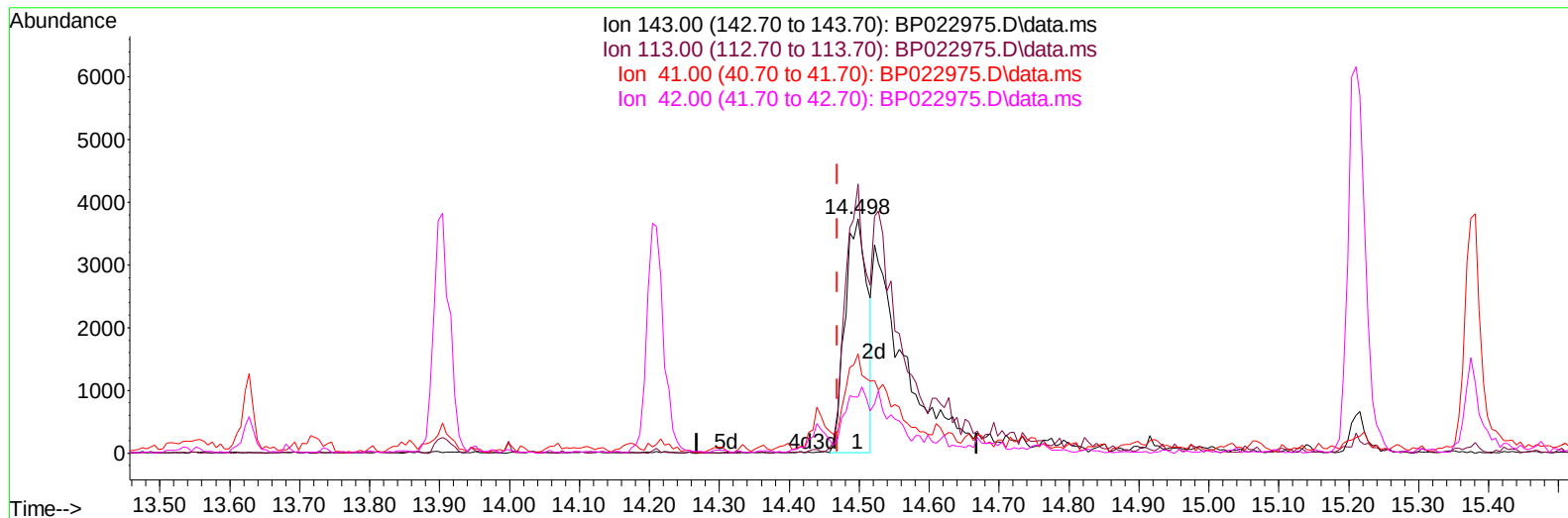
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022975.D
 Acq On : 09 Nov 2024 20:38
 Operator : RC/JU
 Sample : P4746-20
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP93

Manual IntegrationsAPPROVED

Quant Time: Nov 09 23:31:52 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: BP022975.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.029) 4.06 ng/u1

response 8362

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	114.80
41.00	39.20	42.26
42.00	25.90	23.88

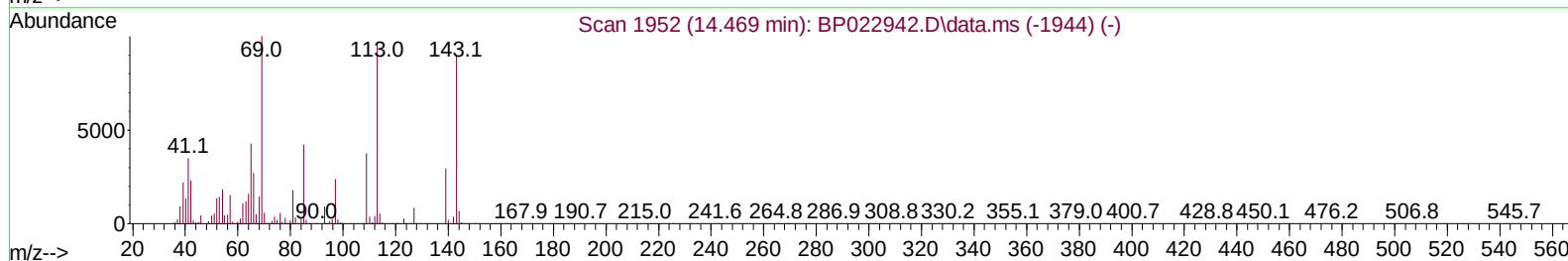
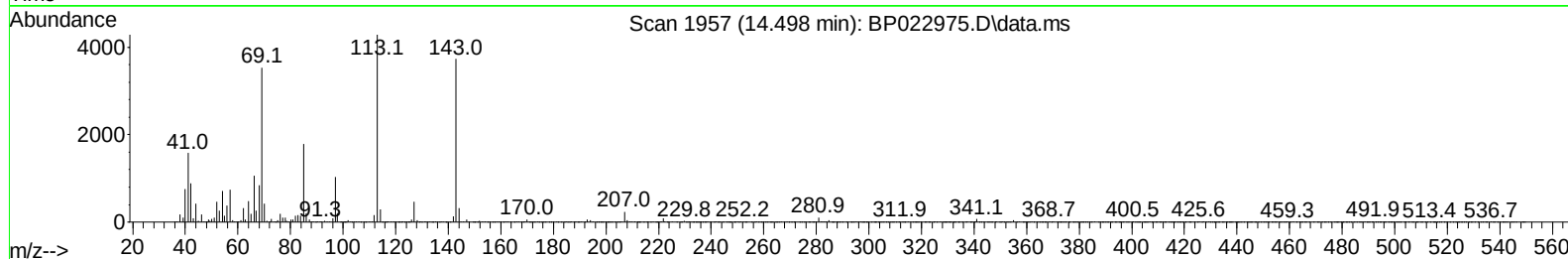
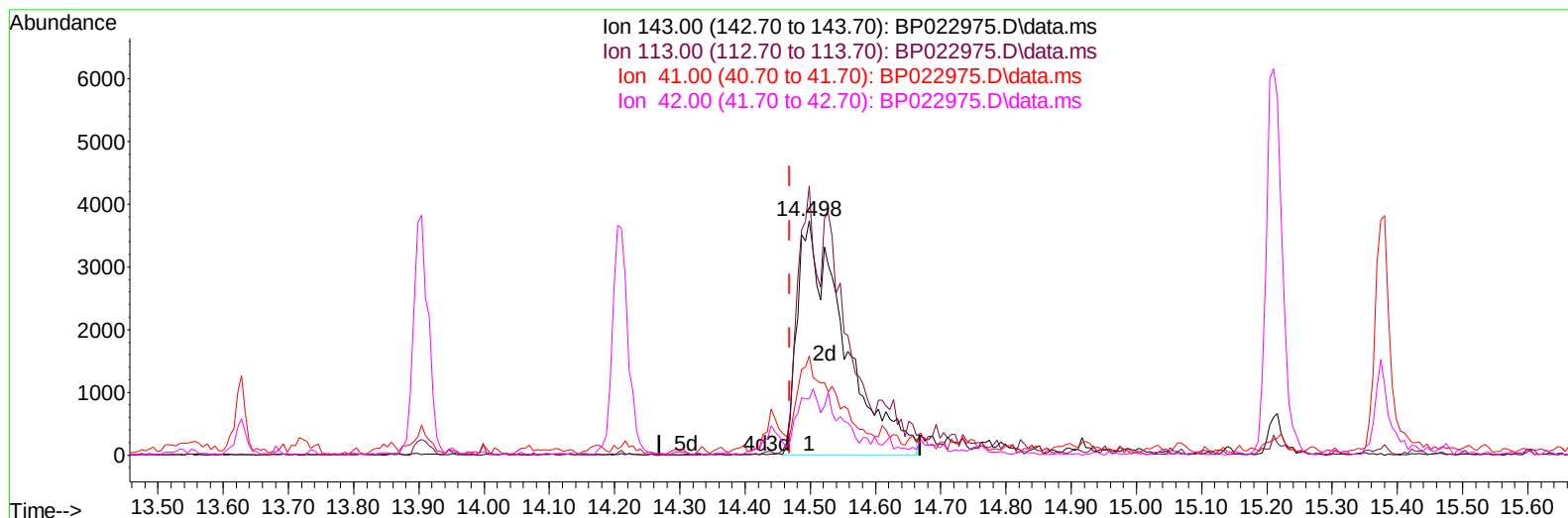
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(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.029) 9.13 ng/ul m

response 18802

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	114.80
41.00	39.20	42.26
42.00	25.90	23.88

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	62226	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	232801	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.210	164	190048	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.004	188	484246	20.000	ng/ul	-0.03
79) Chrysene-d12	21.451	240	604314	20.000	ng/ul	0.00
88) Perylene-d12	24.680	264	713561	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6182	4.775	ng/uL	0.00
4) Pyridine-d5	3.569	84	16603	4.317	ng/ul	0.01
7) Phenol-d5	6.799	99	41079	8.929	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	79723	29.495	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	93405	27.722	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	78178	20.650	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	49927	30.806	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	64348	33.304	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	121745	30.088	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	73674	15.262	ng/ul	0.00
46) Dimethylphthalate-d6	13.628	166	500545	36.543	ng/ul	0.01
49) Acenaphthylene-d8	13.904	160	458444	32.077	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	18802m	9.134	ng/ul	0.03
60) Fluorene-d10	15.210	176	406714	33.867	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.375	200	88686	30.572	ng/ul	0.00
73) Anthracene-d10	17.122	188	739858	36.085	ng/ul	0.00
81) Pyrene-d10	19.498	212	1005917	36.138	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.451	264	1240579	36.499	ng/ul	-0.02

Target Compounds Qvalue

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

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