

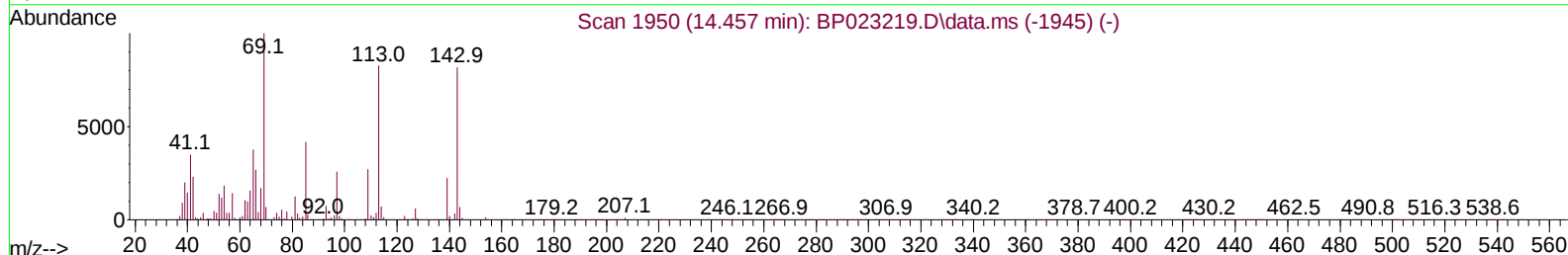
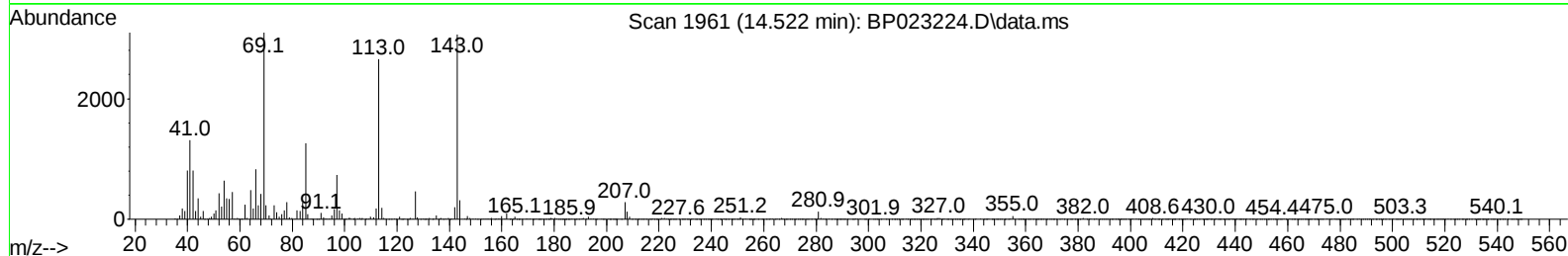
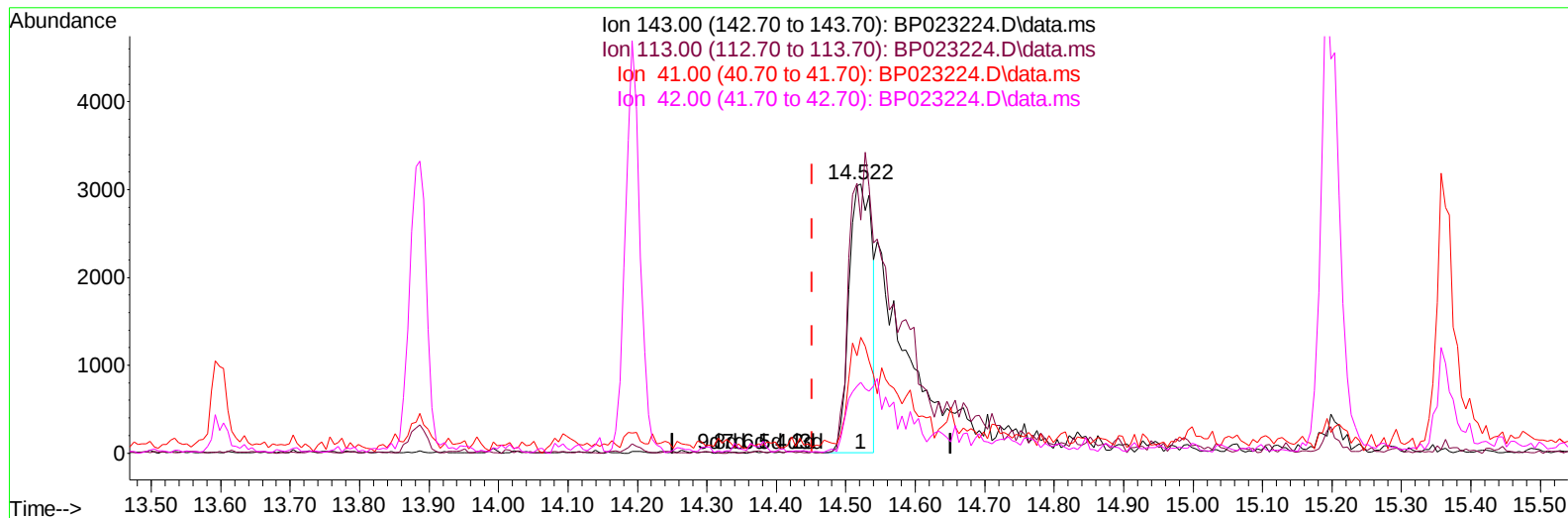
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023224.D
Acq On : 19 Nov 2024 19:21
Operator : RC/JU
Sample : P4875-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29Z5

Manual IntegrationsAPPROVED

Quant Time: Nov 19 23:46:05 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/20/2024
Supervised By :mohammad ahmed 11/22/2024



TIC: BP023224.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.071) 3.07 ng/u1

response 6952

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	86.71
41.00	39.20	42.95
42.00	25.90	26.29

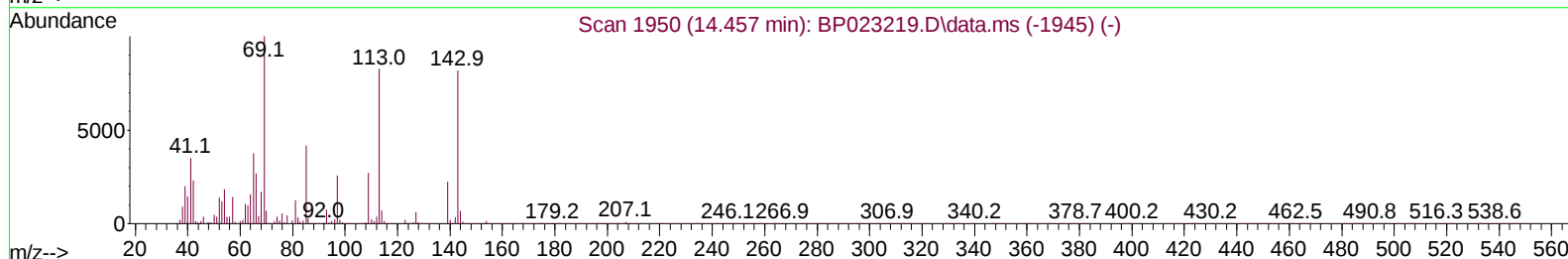
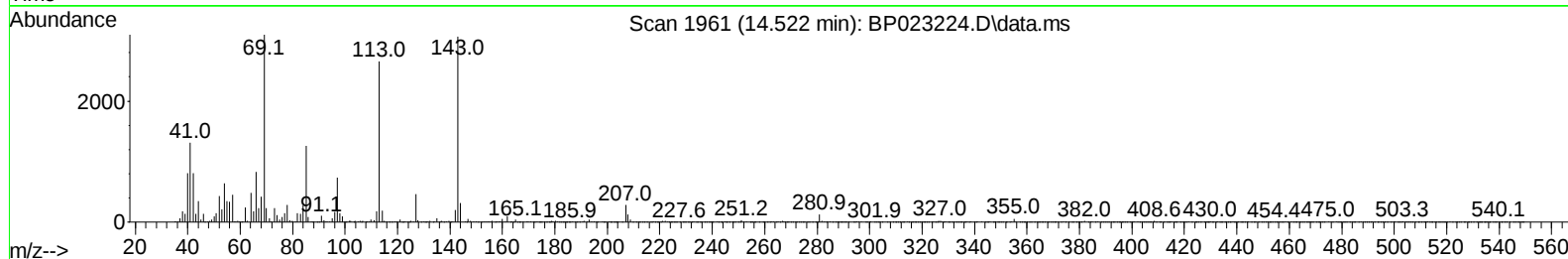
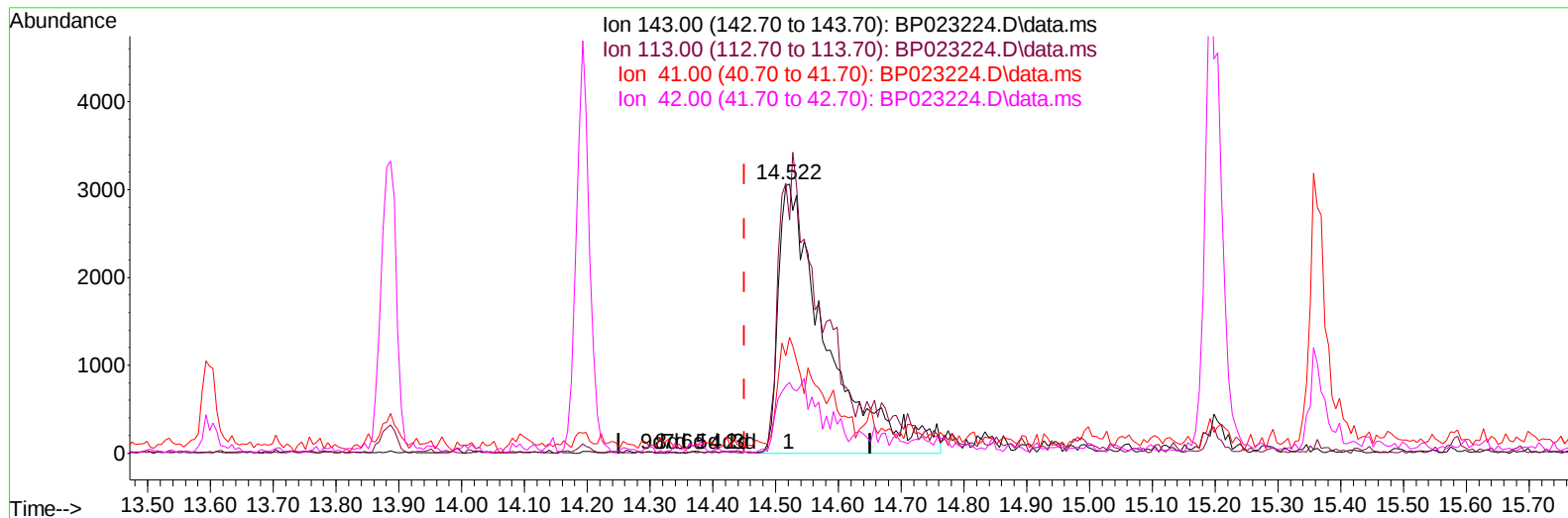
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(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.071) 7.20 ng/ul m

response 16327

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	86.71
41.00	39.20	42.95
42.00	25.90	26.29

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Quant Time: Nov 20 00:06:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	75982	20.000	ng/ul	0.00
20) Naphthalene-d8	10.322	136	270271	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	209251	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.998	188	498685	20.000	ng/ul	-0.02
79) Chrysene-d12	21.439	240	611037	20.000	ng/ul	0.00
88) Perylene-d12	24.651	264	763147	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	7692	4.866	ng/uL	0.00
4) Pyridine-d5	3.575	84	12843	2.735	ng/ul	0.03
7) Phenol-d5	6.793	99	41183	7.331	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.916	67	82608	25.029	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.122	132	92357	22.449	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	78494	16.980	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	47596	25.296	ng/ul	0.00
24) 2-Nitrophenol-d4	9.428	143	55276	24.642	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	114323	24.337	ng/ul	0.01
31) 4-Chloroaniline-d4	10.481	131	85763	15.303	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	447710	29.686	ng/ul	-0.01
49) Acenaphthylene-d8	13.881	160	405685	25.780	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.522	143	16327m	7.204	ng/ul	0.07
60) Fluorene-d10	15.198	176	368028	27.833	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	76441	25.588	ng/ul	0.00
73) Anthracene-d10	17.104	188	696562	32.989	ng/ul	-0.01
81) Pyrene-d10	19.474	212	919771	32.680	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.433	264	1212859	33.365	ng/ul	-0.01

Target Compounds Qvalue

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

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