

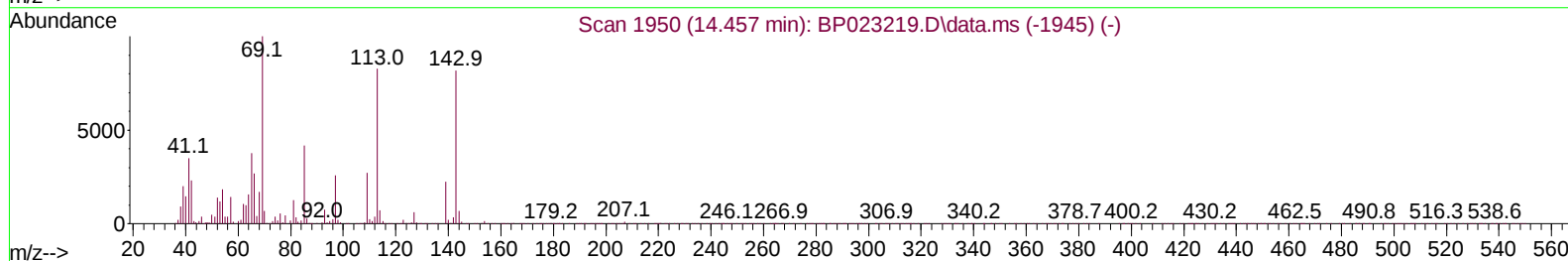
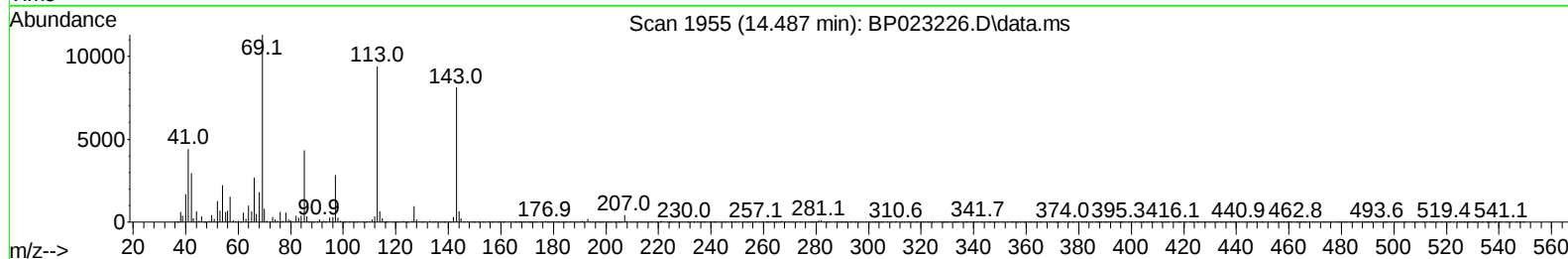
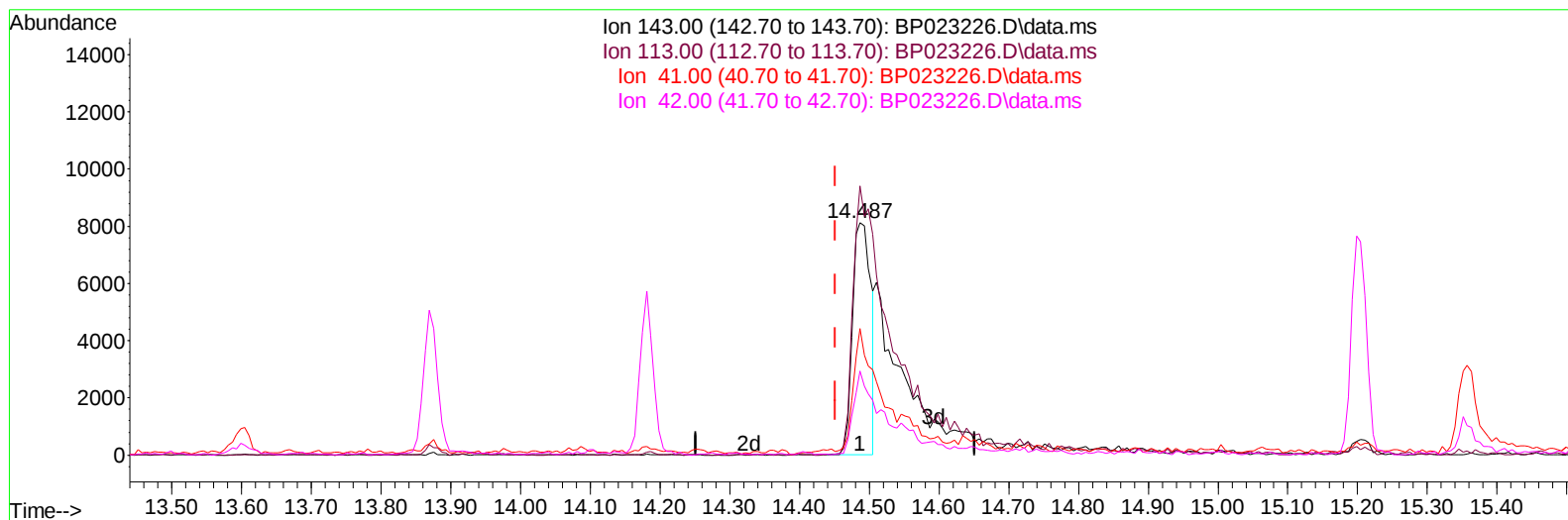
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023226.D
Acq On : 19 Nov 2024 20:42
Operator : RC/JU
Sample : P4875-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E2A13

Manual IntegrationsAPPROVED

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

Quant Time: Nov 20 00:12:14 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



TIC: BP023226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 6.20 ng/u1

response 14712

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	115.84
41.00	39.20	54.50#
42.00	25.90	36.14#

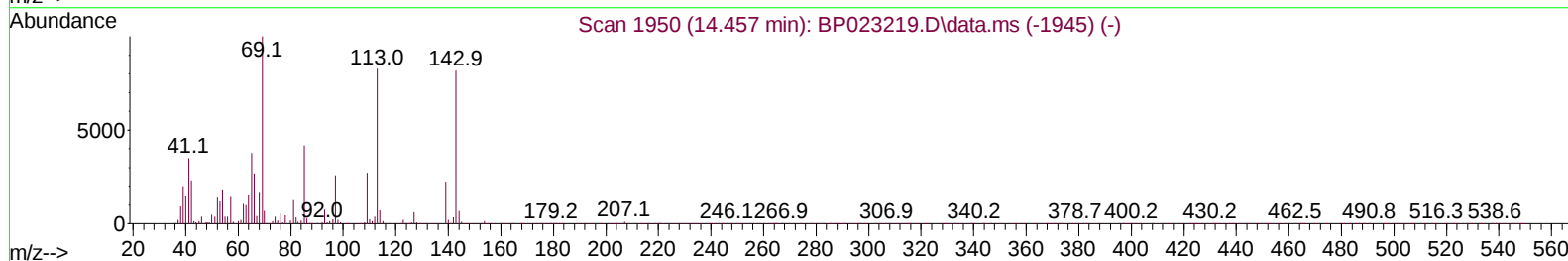
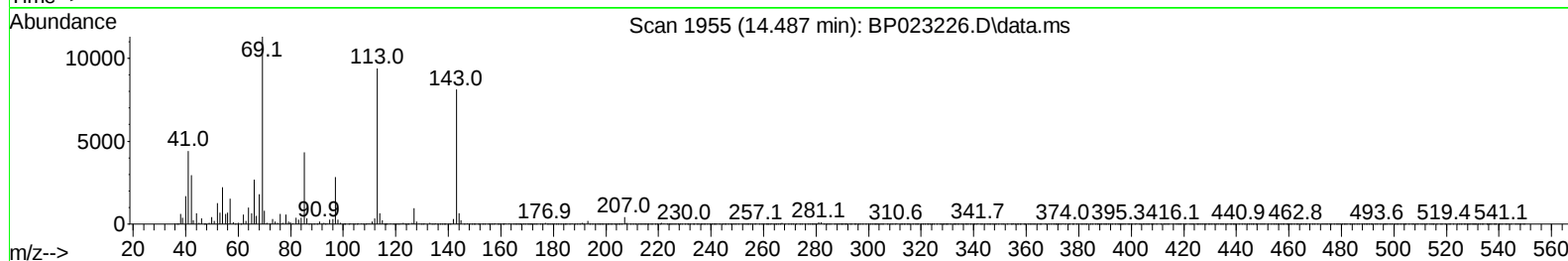
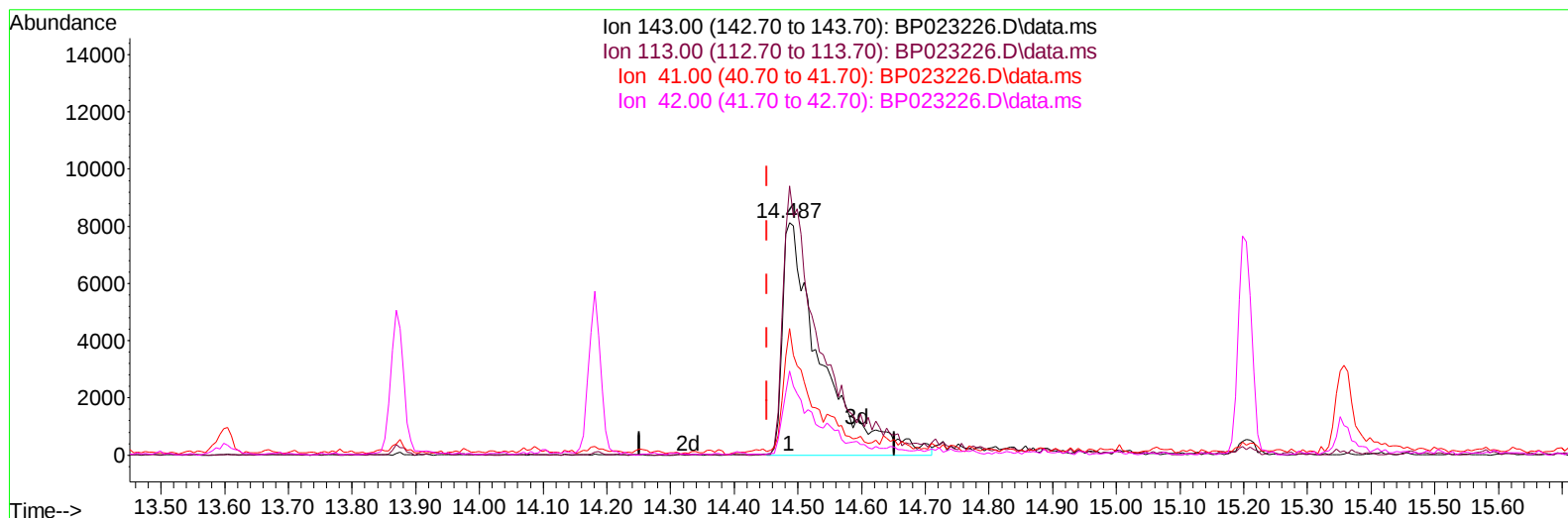
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TIC: BP023226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 14.40 ng/ul m

response 34147

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	115.84
41.00	39.20	54.50#
42.00	25.90	36.14#

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Reviewed By :Yogesh Patel 11/20/2024
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Quant Time: Nov 20 00:13:06 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	80016	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.316	136	287060	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.181	164	218989	20.000	ng/ul	-0.02
64) Phenanthrene-d10	16.998	188	531341	20.000	ng/ul	-0.02
79) Chrysene-d12	21.427	240	652559	20.000	ng/ul	-0.02
88) Perylene-d12	24.639	264	814034	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	9282	5.576	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.781	99	77482	13.097	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.916	67	101636	29.242	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.122	132	127255	29.372	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	115520	23.730	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	60231	30.139	ng/ul	0.00
24) 2-Nitrophenol-d4	9.428	143	72855	30.579	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	145144	29.091	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	68233	11.463	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166	534884	33.889	ng/ul	0.00
49) Acenaphthylene-d8	13.875	160	505311	30.683	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.487	143	34147m	14.396	ng/ul	0.04
60) Fluorene-d10	15.204	176	456920	33.020	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	94172	29.586	ng/ul	0.00
73) Anthracene-d10	17.087	188	809070	35.963	ng/ul	-0.03
81) Pyrene-d10	19.481	212	1089807	36.258	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.416	264	1426466	36.788	ng/ul	-0.03

Target Compounds Qvalue

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

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