

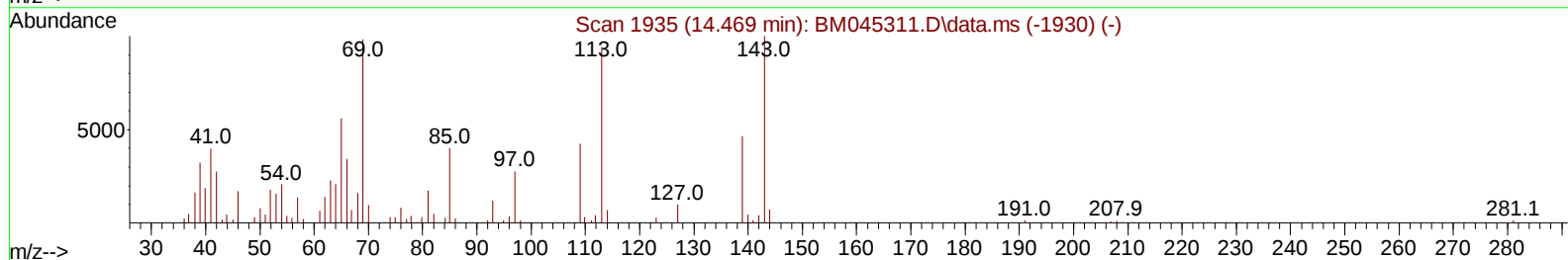
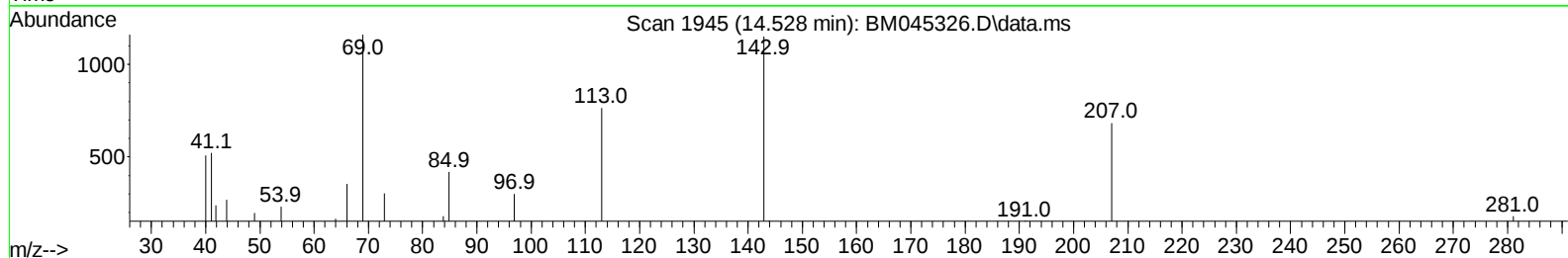
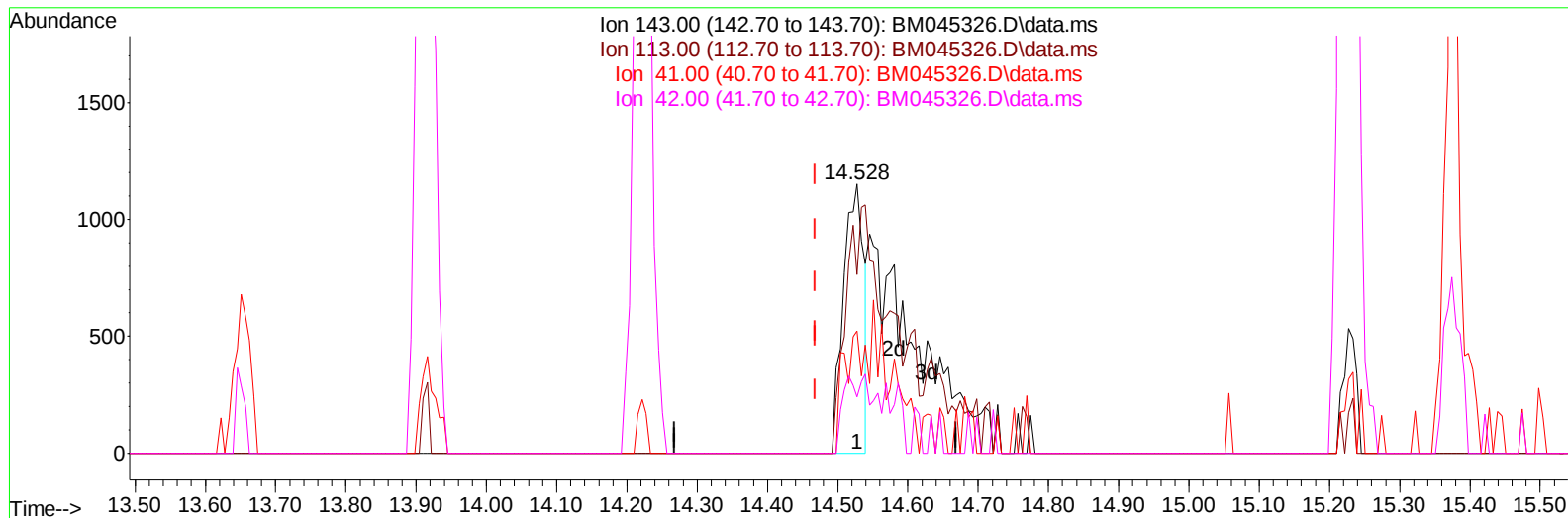
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041724\
Data File : BM045326.D
Acq On : 17 Apr 2024 20:43
Operator : MA/JU
Sample : P2032-22
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0MG6

Manual IntegrationsAPPROVED

Quant Time: Apr 17 23:18:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 17 12:42:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/18/2024
Supervised By :mohammad ahmed 04/25/2024



TIC: BM045326.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.059) 1.21 ng/ul

response 2300

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	66.46
41.00	36.50	45.35#
42.00	23.30	20.85

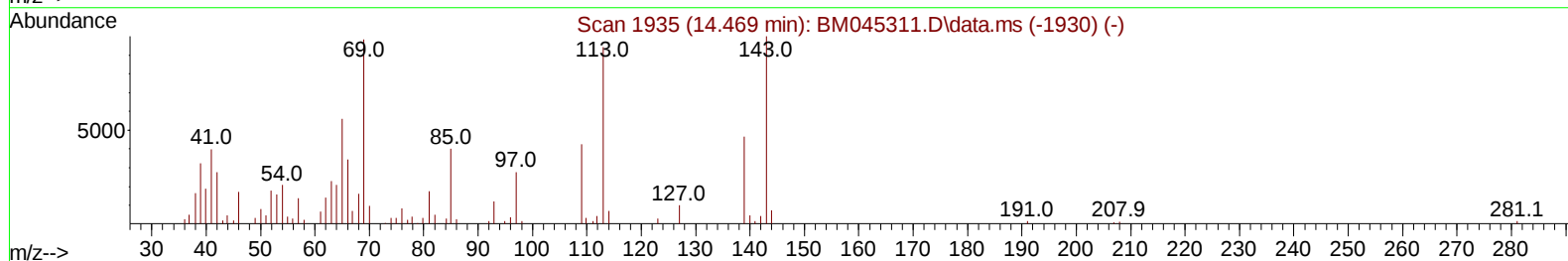
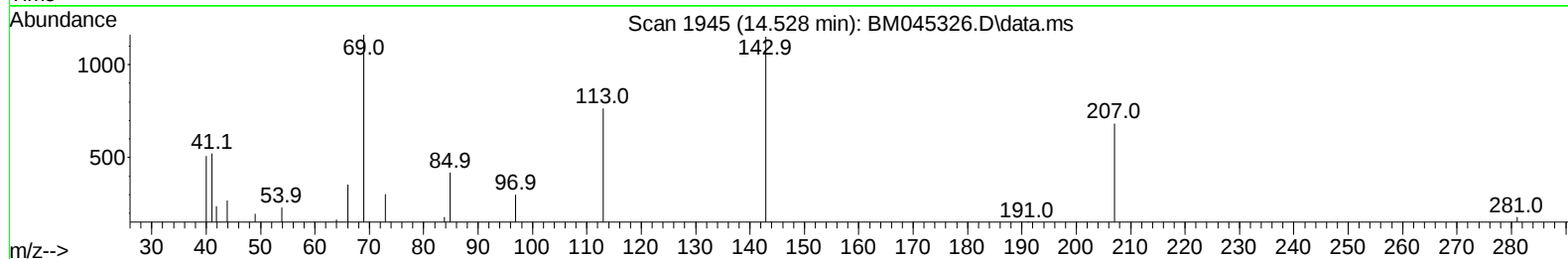
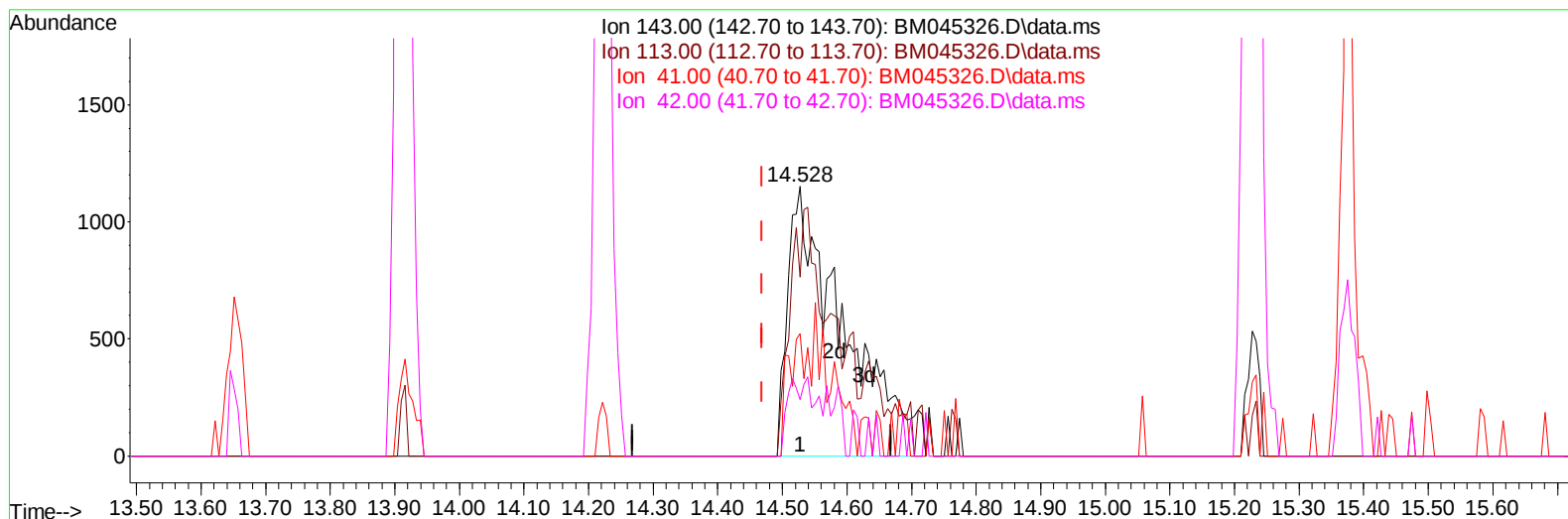
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(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.059) 3.65 ng/ul m

response 6929

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	66.46
41.00	36.50	45.35#
42.00	23.30	20.85

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	62339	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	235104	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.222	164	130779	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.986	188	254119	20.000	ng/ul	0.00
79) Chrysene-d12	21.198	240	243172	20.000	ng/ul	0.00
88) Perylene-d12	23.398	264	344726	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	5671	3.397	ng/uL	0.00
4) Pyridine-d5	3.575	84	37291	8.214	ng/ul	0.00
7) Phenol-d5	6.763	99	32102	6.074	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	90630	28.784	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	96387	23.280	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	57551	13.913	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	56194	31.117	ng/ul	0.00
24) 2-Nitrophenol-d4	9.445	143	57478	30.155	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	84946	24.307	ng/ul	0.00
31) 4-Chloroaniline-d4	10.504	131	121947	21.917	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	295893	32.495	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	337567	31.380	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	6929m	3.646	ng/ul	0.06
60) Fluorene-d10	15.227	176	253050	31.171	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.374	200	33251	22.068	ng/ul	0.00
73) Anthracene-d10	17.086	188	389630	34.084	ng/ul	0.00
81) Pyrene-d10	19.392	212	454844	37.987	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.256	264	588849	35.016	ng/ul	0.00

Target Compounds Qvalue

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

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