

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041524\
Data File : BM045246.D
Acq On : 15 Apr 2024 16:13
Operator : MA/JU
Sample : PB160235BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

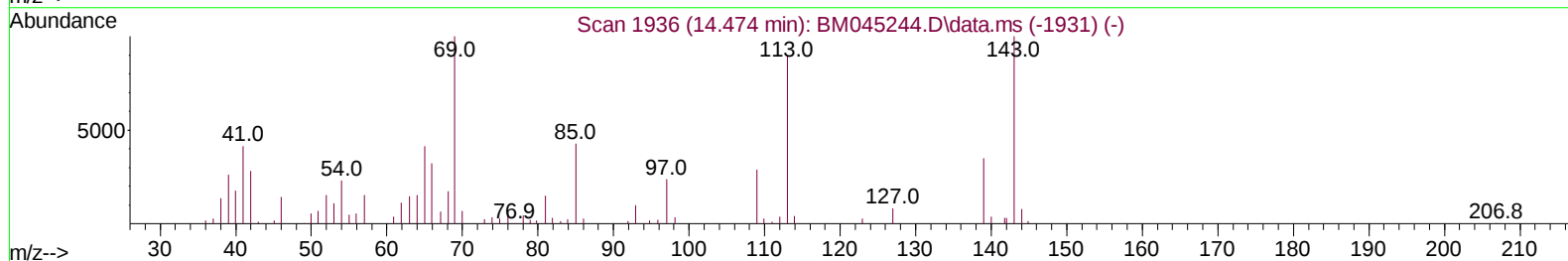
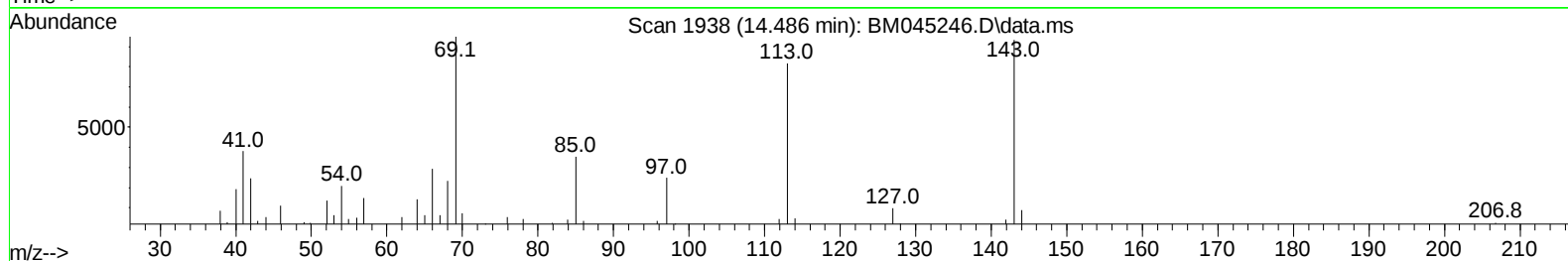
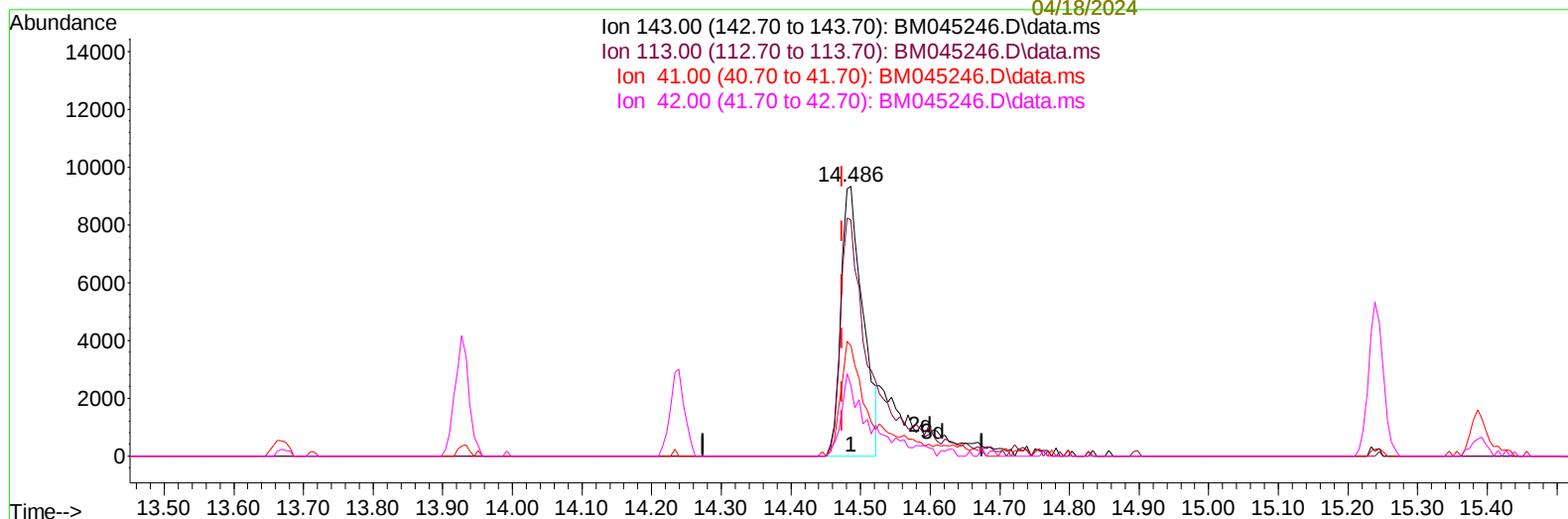
Instrument :
BNA_M
ClientSampleId :
SBLK235

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/17/2024
Supervised By :mohammad ahmed 04/18/2024

Supervised By :mohammad
ahmed
04/18/2024

Quant Time: Apr 15 17:18:32 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 15 16:06:37 2024
Response via : Initial Calibration



TIC: BM045246.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.012) 18.87 ng/u1

response 20406

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	87.60
41.00	36.50	40.88
42.00	23.30	26.28

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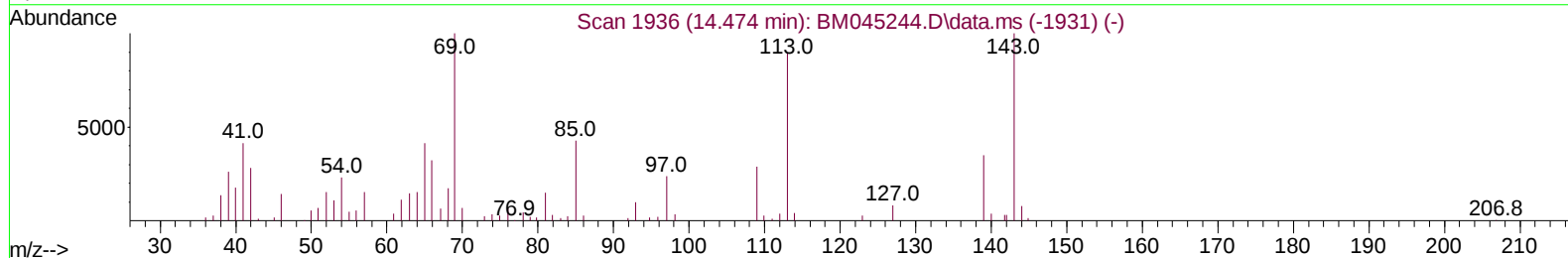
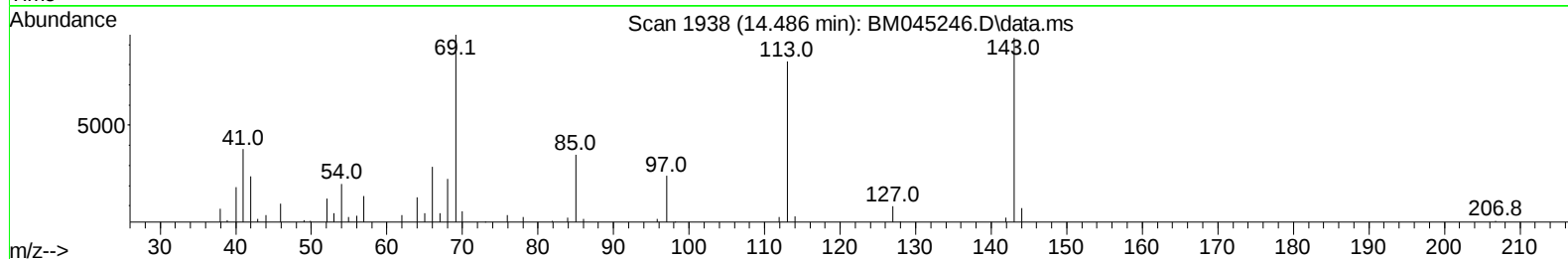
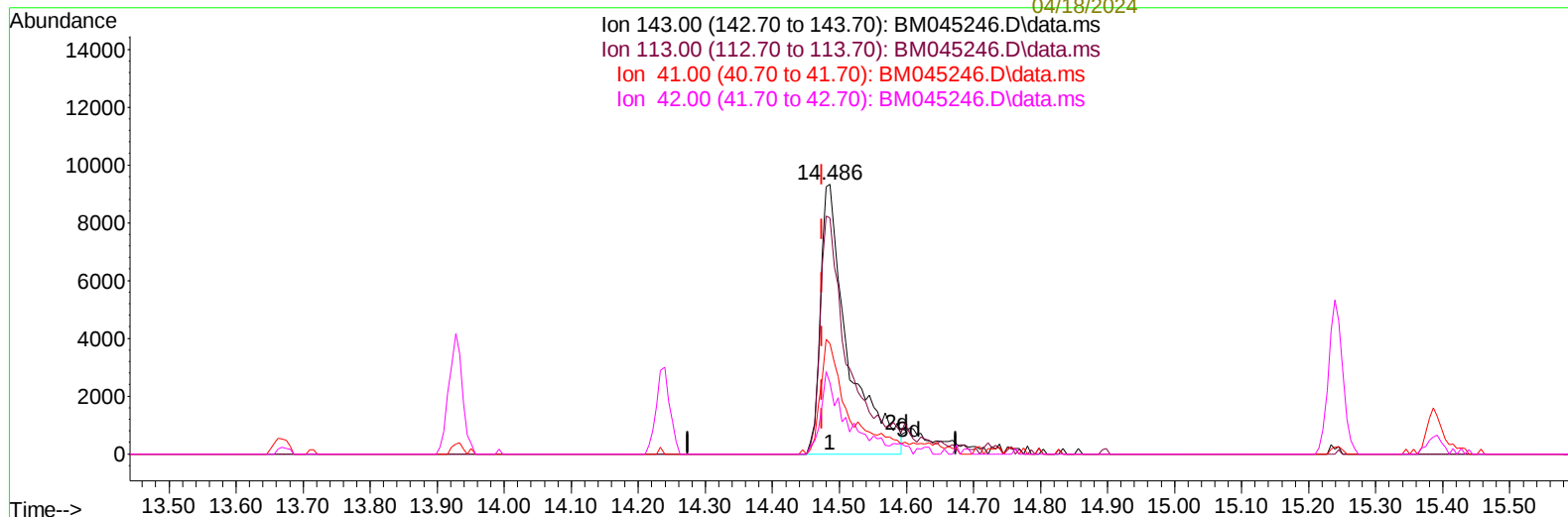
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TIC: BM045246.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.012) 24.63 ng/ul m

response 26639

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	87.60
41.00	36.50	40.88
42.00	23.30	26.28

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Quant Time: Apr 15 17:19:09 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.587	152		34502	20.000	ng/ul	0.00
20) Naphthalene-d8	10.357	136		131732	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.239	164		74431	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188		150317	20.000	ng/ul	0.00
79) Chrysene-d12	21.215	240		140786	20.000	ng/ul	0.00
88) Perylene-d12	23.421	264		187358	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.181	96		6398	6.925	ng/uL	0.00
4) Pyridine-d5	3.575	84		84387	33.586	ng/ul	0.00
7) Phenol-d5	6.769	99		100739	34.442	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67		65844	37.784	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132		88669	38.695	ng/ul	0.00
15) 4-Methylphenol-d8	8.298	113		78502	34.289	ng/ul	0.00
21) Nitrobenzene-d5	8.751	128		39909	39.441	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143		41669	39.015	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.992	165		65871	33.639	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131		103883	33.321	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166		215878	41.656	ng/ul	0.00
49) Acenaphthylene-d8	13.927	160		242702	39.642	ng/ul	0.00
54) 4-Nitrophenol-d4	14.486	143		26639m	24.630	ng/ul	0.01
60) Fluorene-d10	15.239	176		184765	39.990	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200		24245	27.203	ng/ul	0.00
73) Anthracene-d10	17.098	188		280414	41.469	ng/ul	0.00
81) Pyrene-d10	19.404	212		318927	46.006	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.274	264		386307	42.266	ng/ul	0.00

Target Compounds Qvalue

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

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