

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047764.D
Acq On : 02 Oct 2024 06:12
Operator : RC/JU
Sample : PB163513BL
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

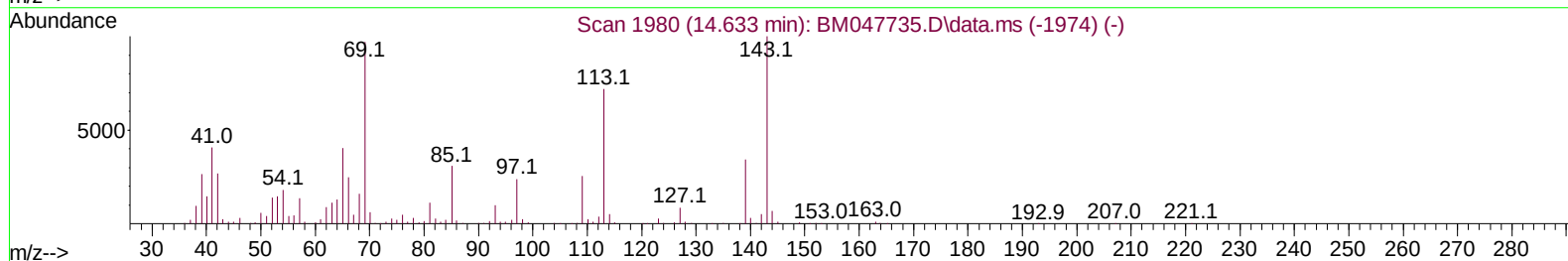
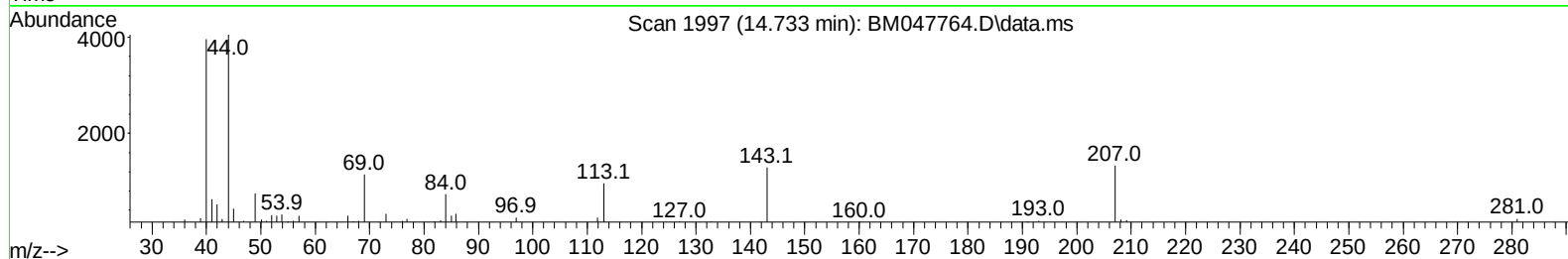
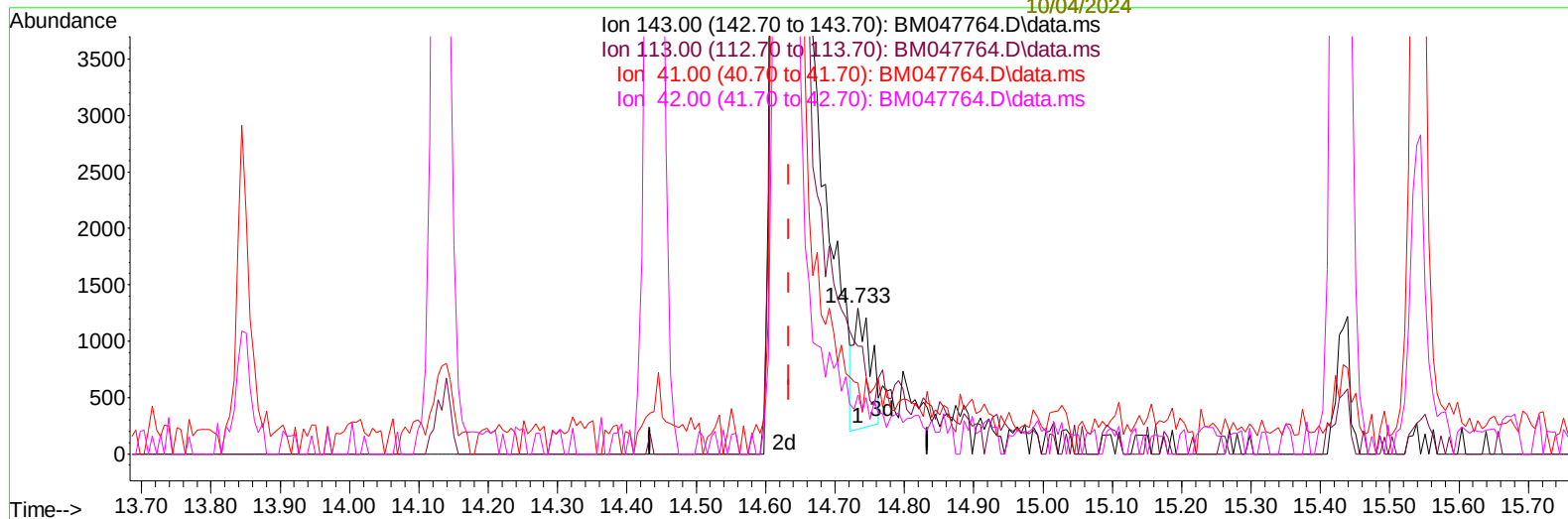
SBLK513

Manual IntegrationsAPPROVED

Quant Time: Oct 02 06:54:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 28 02:30:36 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/04/2024
Supervised By :mohammad ahmed 10/04/2024

Supervised By :mohammad
ahmed
10/04/2024



TIC: BM047764.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.100) 0.21 ng/ul

response 1751

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.20	73.78
41.00	60.00	48.41
42.00	40.50	40.53

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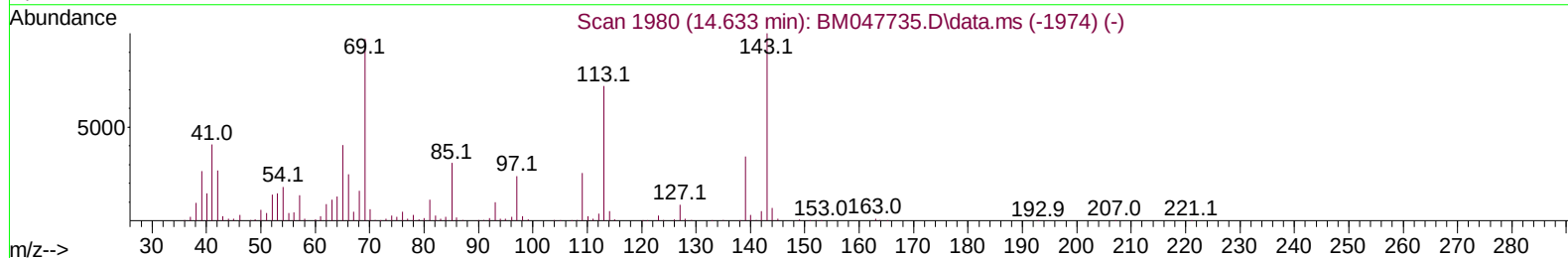
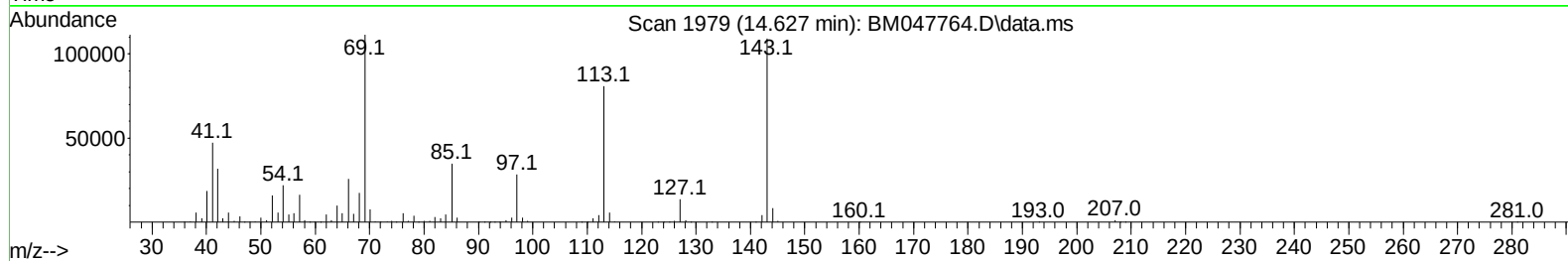
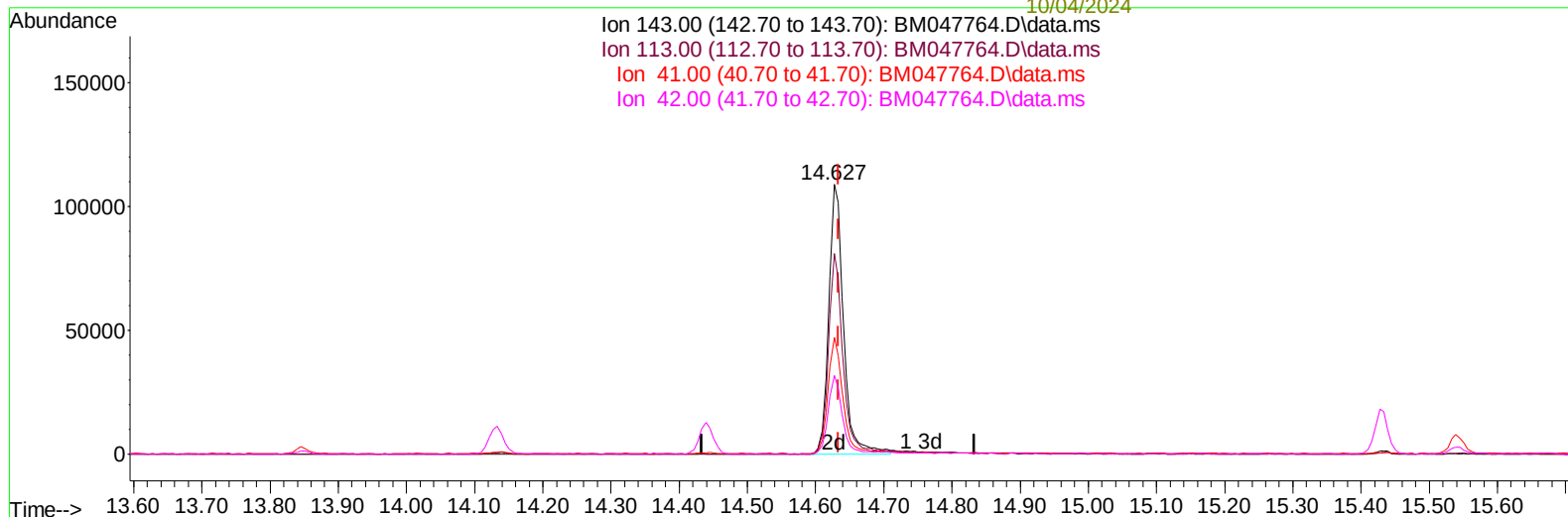
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(54) 4-Nitrophenol-d4 (S)

14.627min (-0.006) 19.40 ng/ul m

response 162634

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.20	74.30
41.00	60.00	43.31#
42.00	40.50	29.18#

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Quant Time: Oct 02 06:55:17 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.810	152		234143	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136		959873	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.439	164		580684	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.180	188		1166614	20.000	ng/ul	0.00
79) Chrysene-d12	21.403	240		1069806	20.000	ng/ul	0.00
88) Perylene-d12	24.403	264		1250797	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.257	96		33178	4.579	ng/uL	0.00
4) Pyridine-d5	3.669	84		469938	22.152	ng/ul	0.00
7) Phenol-d5	6.975	99		506575	22.880	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67		319695	20.348	ng/ul	0.00
11) 2-Chlorophenol-d4	7.340	132		421369	26.158	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113		388338	23.591	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128		195161	25.536	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143		195568	26.703	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.222	165		365609	24.990	ng/ul	0.00
31) 4-Chloroaniline-d4	10.733	131		554362	24.509	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166		1054209	25.041	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160		1251438	25.071	ng/ul	0.00
54) 4-Nitrophenol-d4	14.627	143		162634m	19.403	ng/ul	0.00
60) Fluorene-d10	15.433	176		919538	25.216	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200		107925	15.809	ng/ul	0.00
73) Anthracene-d10	17.274	188		1387844	24.194	ng/ul	-0.01
81) Pyrene-d10	19.562	212		1575947	25.956	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.203	264		1579971	24.346	ng/ul	0.00

Target Compounds Qvalue

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

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