

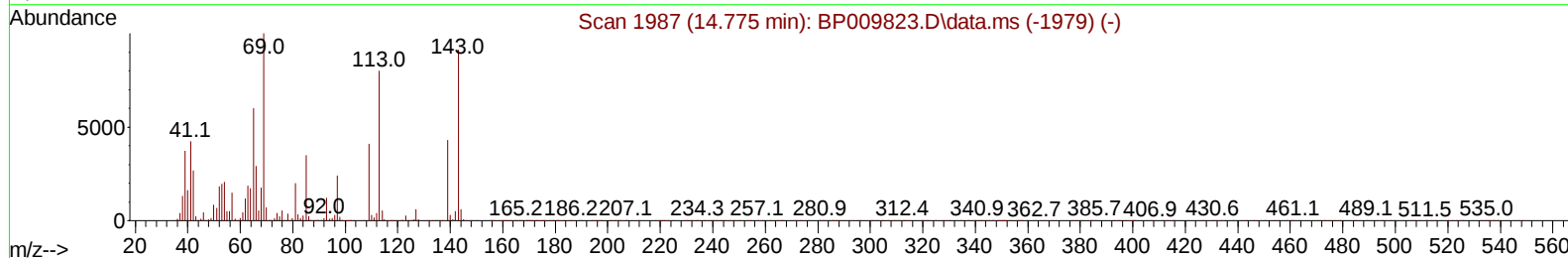
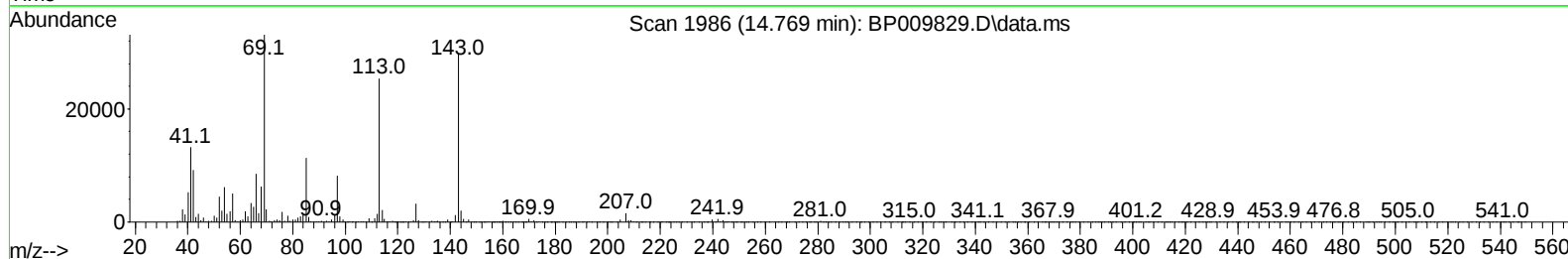
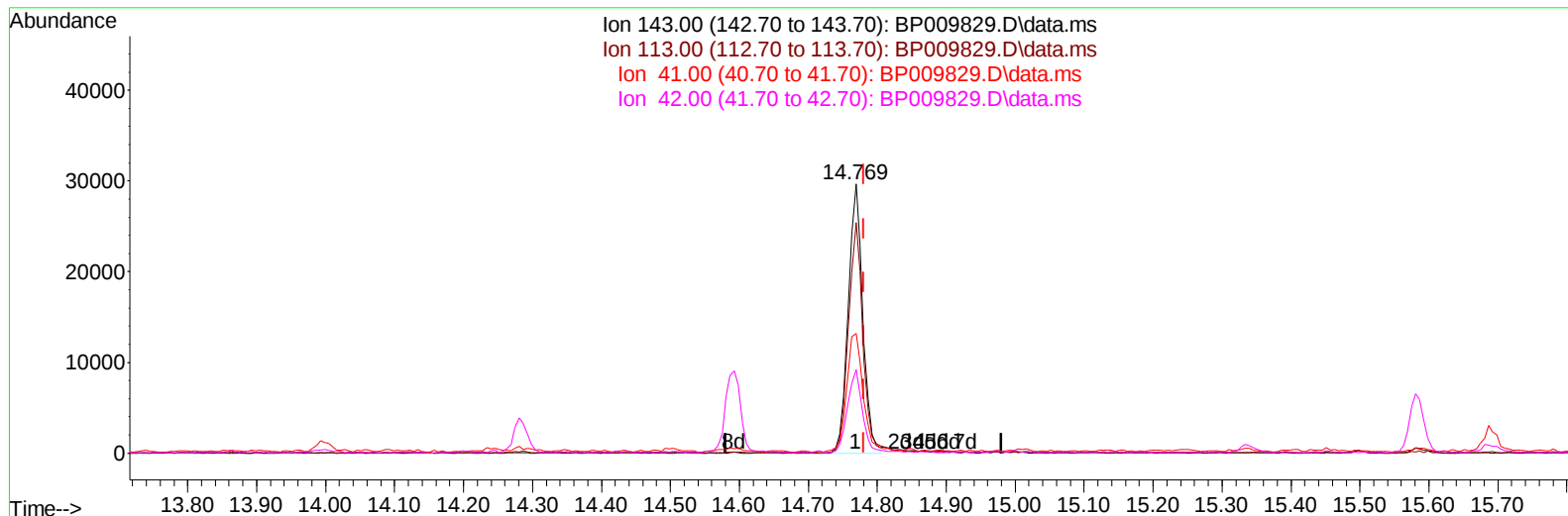
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/22/2022

Quant Time: Apr 14 21:46:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration



TIC: BP009829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 7.20 ng/u1

response 43278

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.67#
41.00	23.20	44.52#
42.00	18.00	31.02#

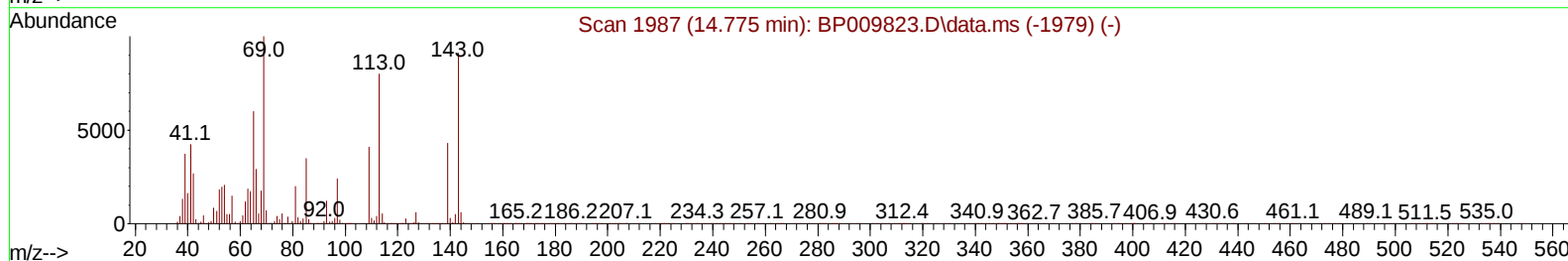
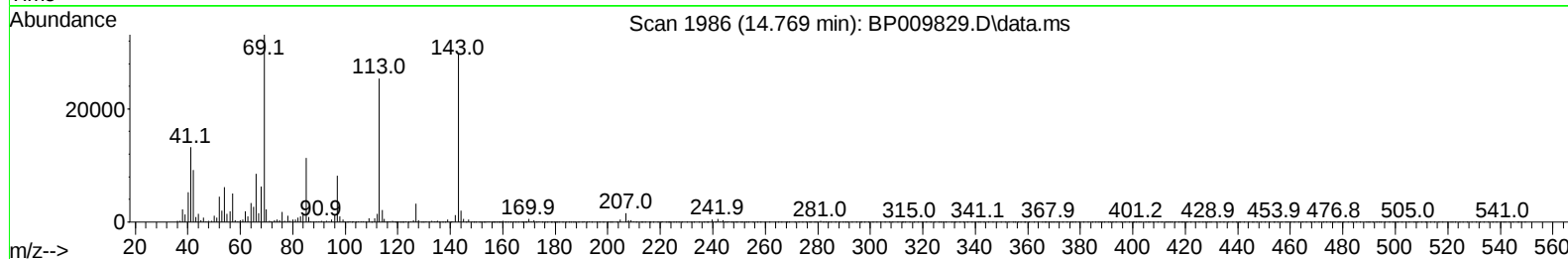
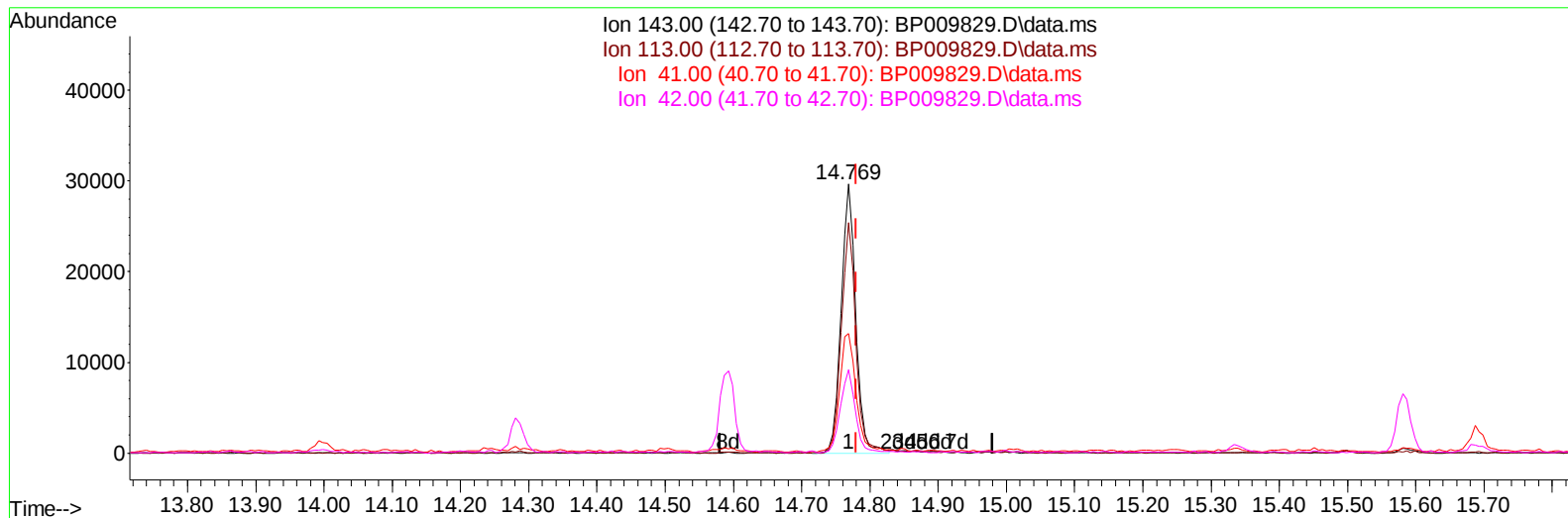
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TIC: BP009829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 7.31 ng/ul m

response 43920

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.67#
41.00	23.20	44.52#
42.00	18.00	31.02#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	133936	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.769	136	591035	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	421736	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	956790	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	993270	20.000	ng/ul	-0.01
88) Perylene-d12	23.886	264	970715	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	5714	1.904	ng/uL	0.00
4) Pyridine-d5	3.805	84	38135	4.609	ng/ul	0.00
7) Phenol-d5	7.116	99	114312	9.721	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	76840	10.977	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	91886	10.359	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	101587	10.349	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	47802	11.216	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	48975	10.481	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	94016	9.615	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.899	131	138853	9.970	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	409610	12.365	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	408014	10.396	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	43920m	7.310	ng/ul	-0.01
60) Fluorene-d10	15.581	176	346448	12.432	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	39319	6.937	ng/ul	0.00
73) Anthracene-d10	17.439	188	624945	13.642	ng/ul	-0.01
81) Pyrene-d10	19.669	212	817518	15.070	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	743701	14.712	ng/ul	-0.01

Target Compounds Qvalue

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

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