

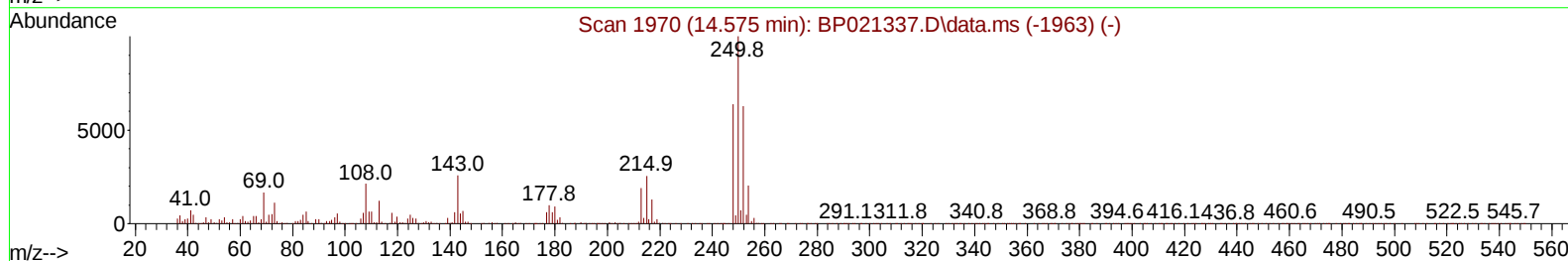
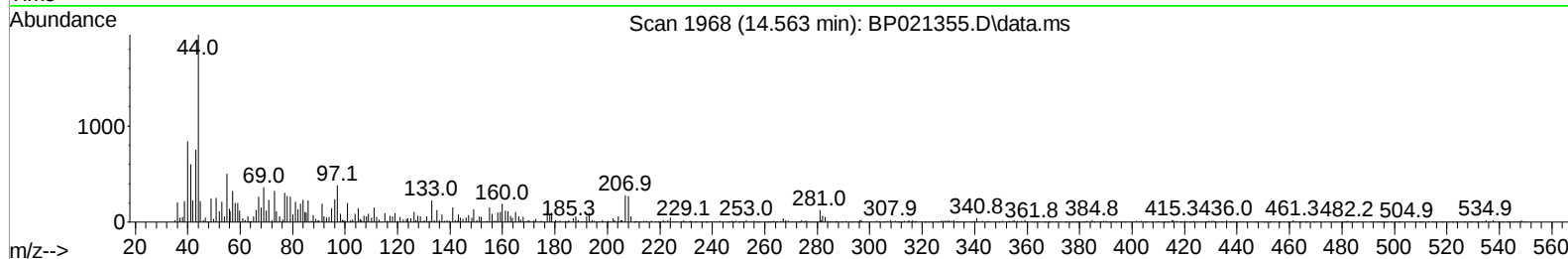
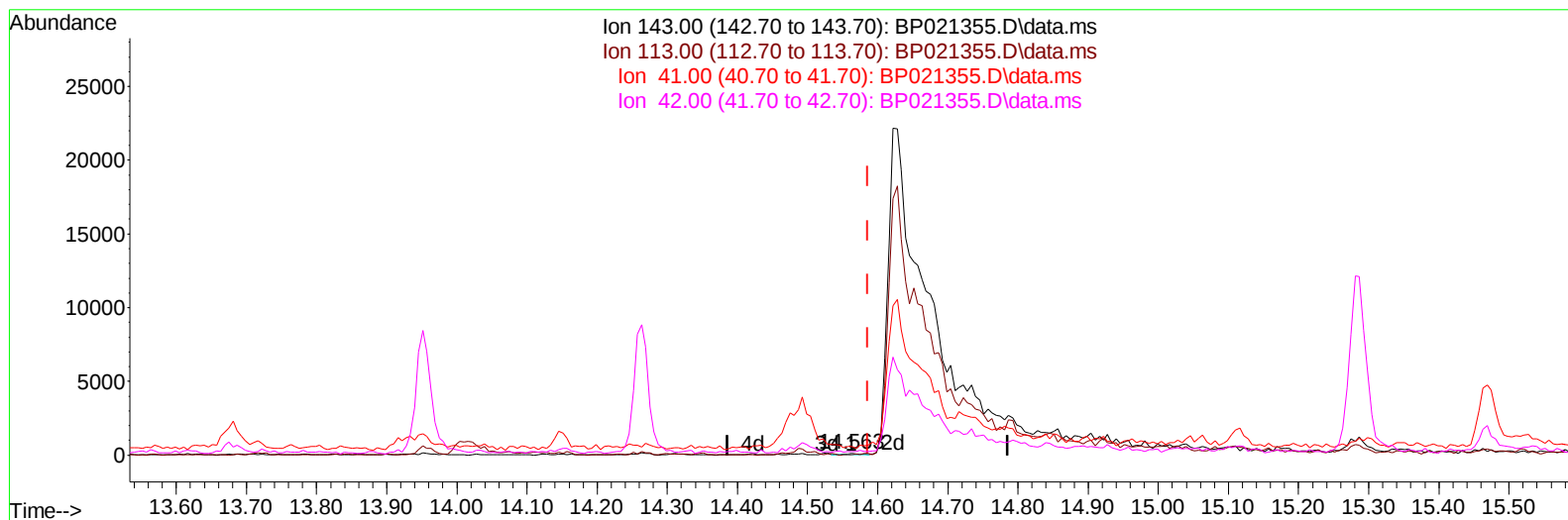
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080324\
Data File : BP021355.D
Acq On : 03 Aug 2024 23:42
Operator : CG/JU
Sample : P3278-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1ZQ2

Manual IntegrationsAPPROVED

Quant Time: Aug 04 23:24:47 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/05/2024
Supervised By :mohammad ahmed 08/05/2024



TIC: BP021355.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (-0.023) 0.03 ng/u1

response 138

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	29.90#
41.00	37.90	619.59#
42.00	27.10	230.93#

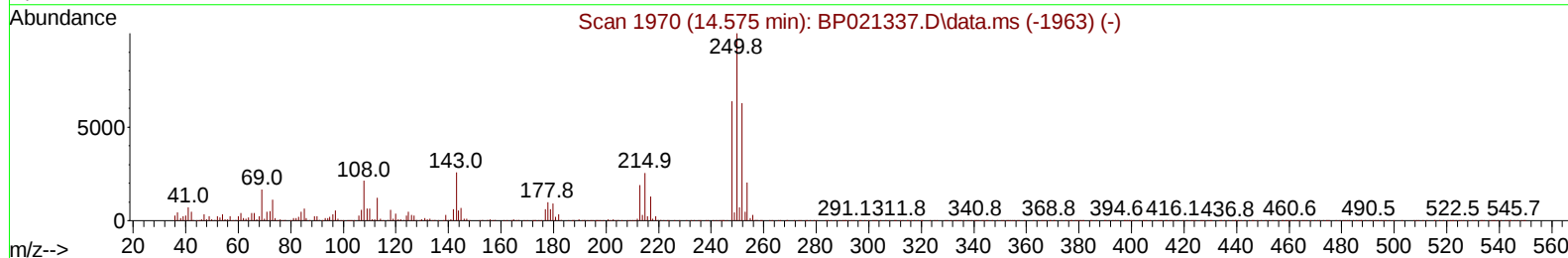
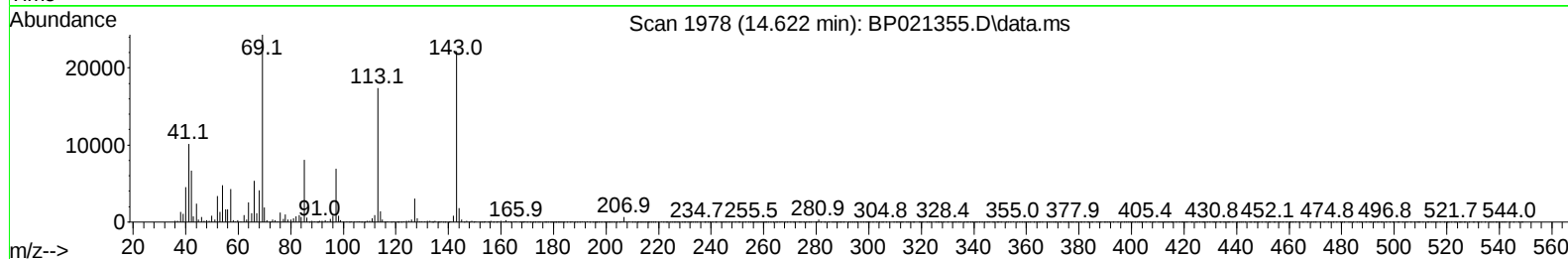
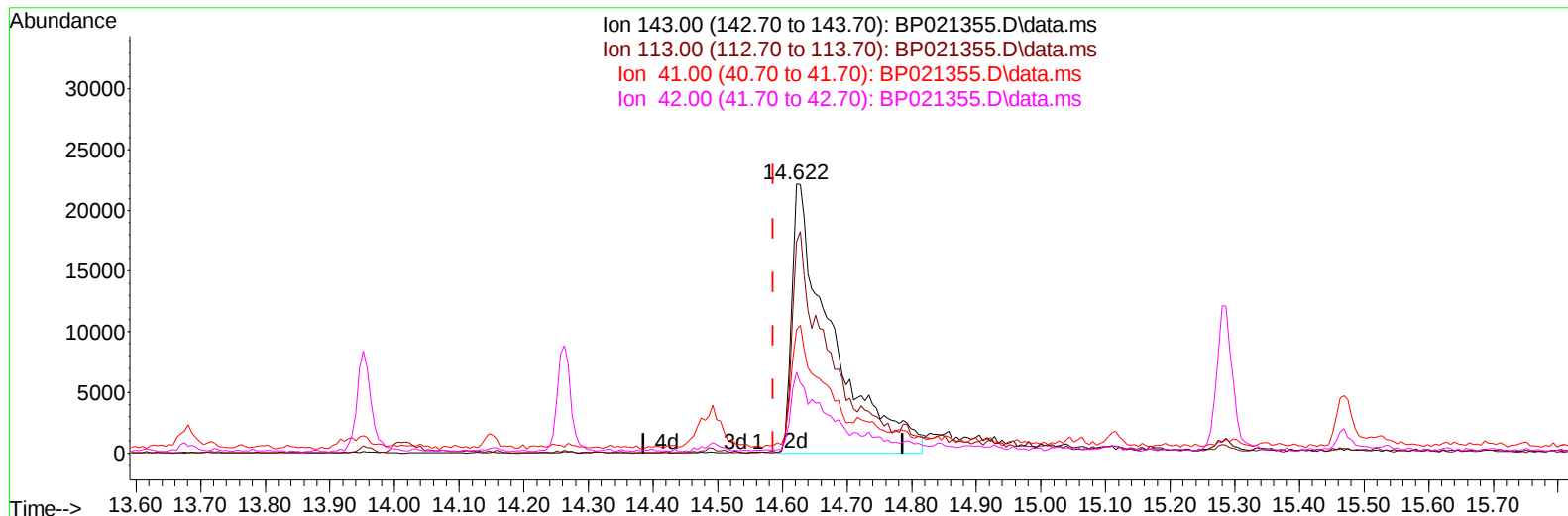
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(54) 4-Nitrophenol-d4 (S)

14.622min (+ 0.036) 22.39 ng/ul m

response 95390

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	78.52
41.00	37.90	45.76#
42.00	27.10	30.04

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Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/05/2024
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Quant Time: Aug 05 00:22:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	184985	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.387	136	690200	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	411988	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	893272	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	1019116	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264	1219014	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	22284	5.482	ng/uL	-0.01
4) Pyridine-d5	3.581	84	263151	22.360	ng/ul	-0.01
7) Phenol-d5	6.846	99	380895	25.256	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	244613	26.764	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	331841	26.706	ng/ul	-0.01
15) 4-Methylphenol-d8	8.358	113	316828	25.294	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	153573	28.647	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	171394	27.933	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	287973	25.955	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	371804	25.296	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	923046	30.820	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1037460	30.850	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.622	143	95390m	22.386	ng/ul	0.04
60) Fluorene-d10	15.287	176	801684	32.628	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	89977	16.648	ng/ul	0.00
73) Anthracene-d10	17.186	188	1264378	31.869	ng/ul	0.00
81) Pyrene-d10	19.569	212	1772093	28.126	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	2073672	34.491	ng/ul	-0.01

Target Compounds Qvalue

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

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