

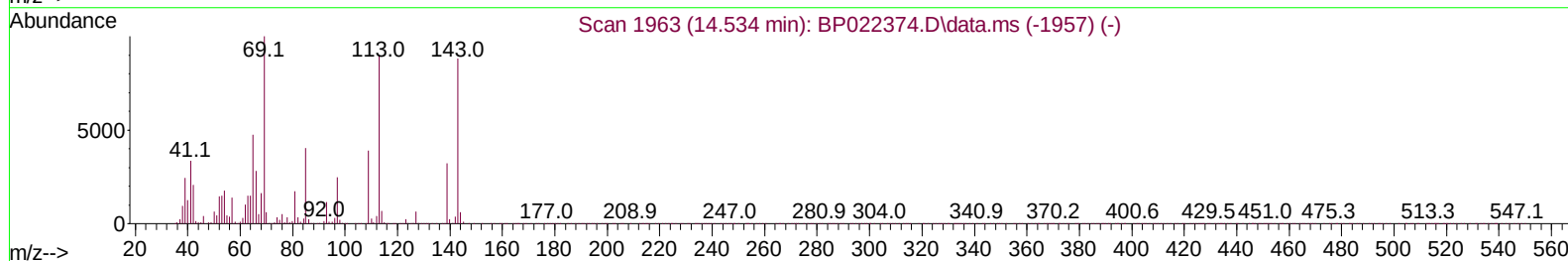
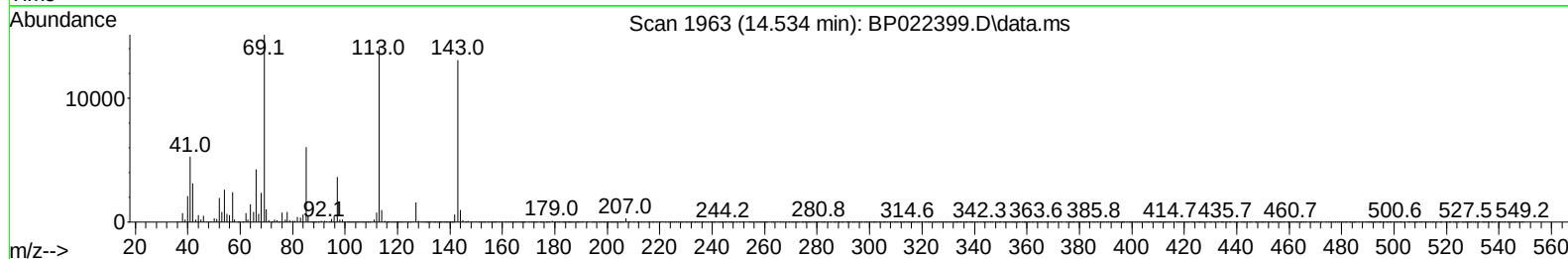
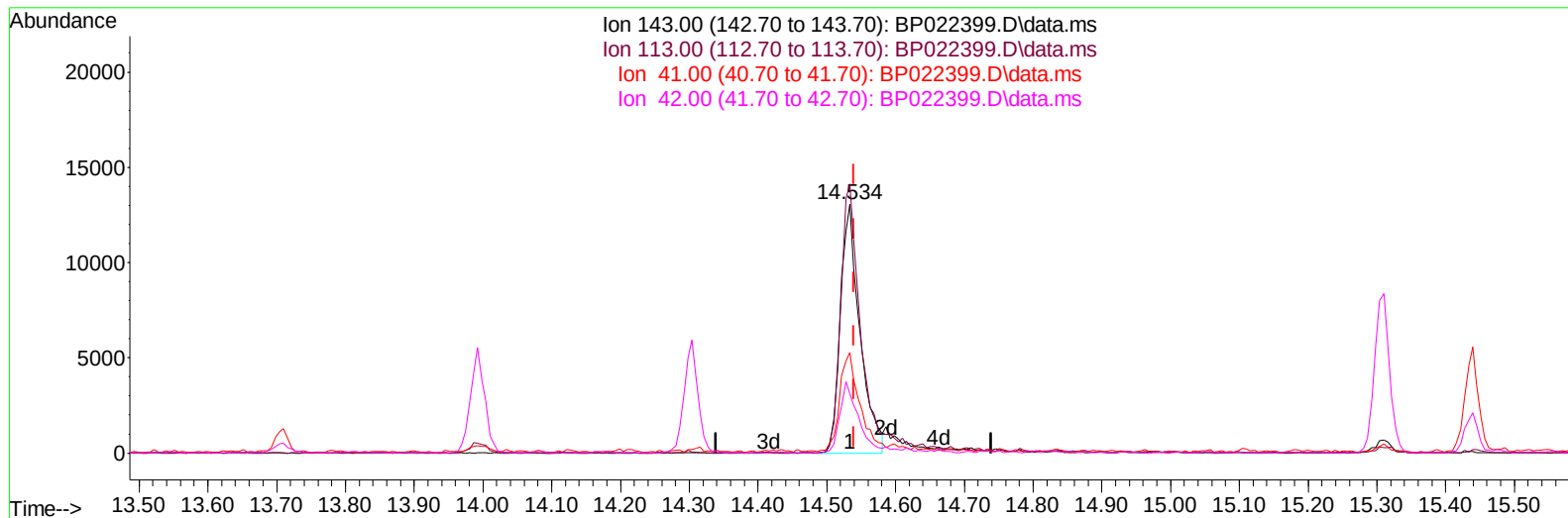
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022399.D
Acq On : 15 Oct 2024 21:42
Operator : RC/JU
Sample : P4349-18
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EOAG3

Manual IntegrationsAPPROVED

Quant Time: Oct 16 02:45:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 14 11:01:55 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/16/2024
Supervised By :mohammad ahmed 10/22/2024



TIC: BP022399.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.534min (-0.006) 7.36 ng/ul

response 25851

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	108.05
41.00	35.90	40.39
42.00	23.70	23.75

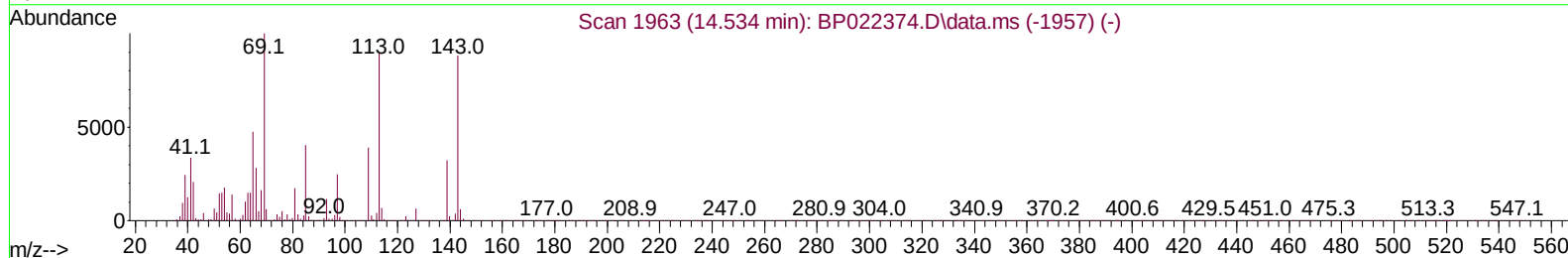
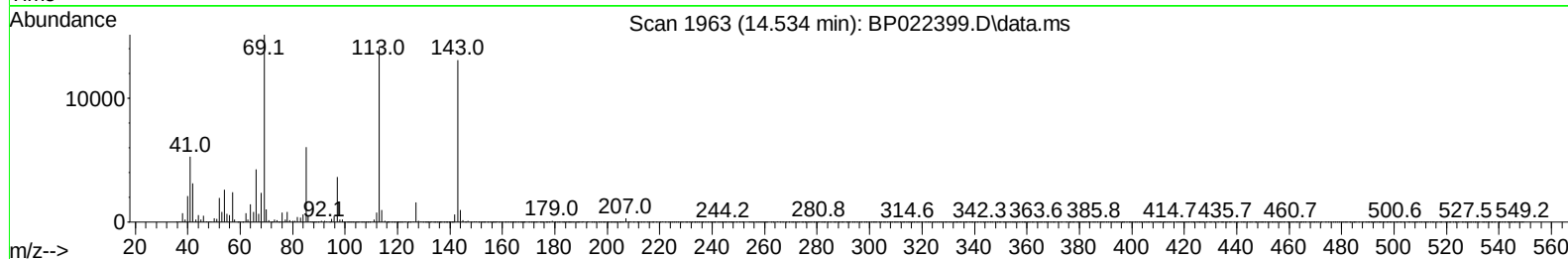
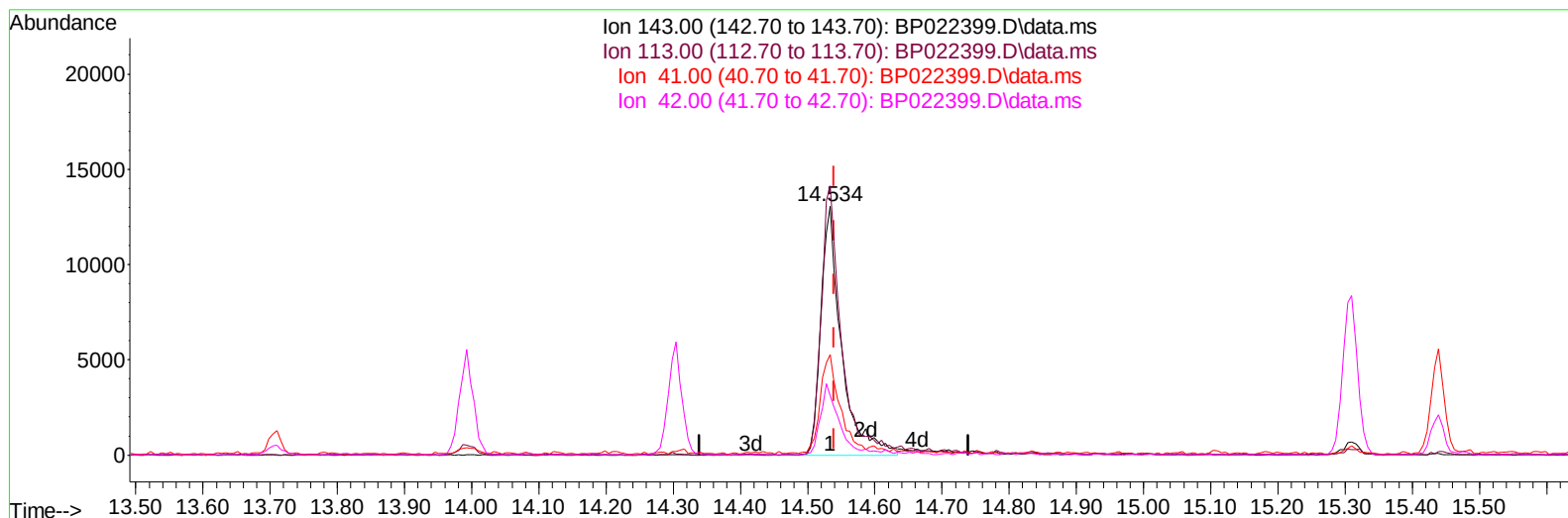
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TIC: BP022399.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.534min (-0.006) 7.96 ng/ul m

response 27926

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	108.05
41.00	35.90	40.39
42.00	23.70	23.75

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Quant Time: Oct 16 02:46:04 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	83284	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	325651	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.304	164	263333	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	669981	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	743192	20.000	ng/ul	-0.01
88) Perylene-d12	24.821	264	868443	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	8393	3.916	ng/uL	0.00
4) Pyridine-d5	3.646	84	25264	4.294	ng/ul	0.01
7) Phenol-d5	6.887	99	56085	8.000	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	113312	27.500	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	131374	26.409	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	101906	18.214	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	72589	28.990	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143	89218	30.776	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	182318	29.591	ng/ul	0.00
31) 4-Chloroaniline-d4	10.598	131	46348	6.366	ng/ul	0.01
46) Dimethylphthalate-d6	13.710	166	631795	30.197	ng/ul	0.00
49) Acenaphthylene-d8	13.992	160	651832	29.333	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	27926m	7.956	ng/ul	0.00
60) Fluorene-d10	15.310	176	570736	31.450	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.439	200	123195	27.958	ng/ul	0.00
73) Anthracene-d10	17.198	188	1053283	33.419	ng/ul	0.00
81) Pyrene-d10	19.574	212	1324500	33.701	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.604	264	1593276	34.353	ng/ul	-0.03

Target Compounds Qvalue

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

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