

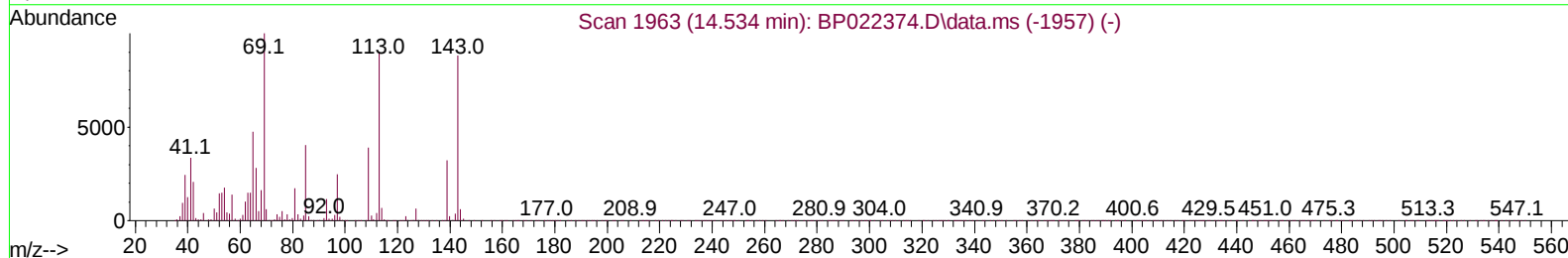
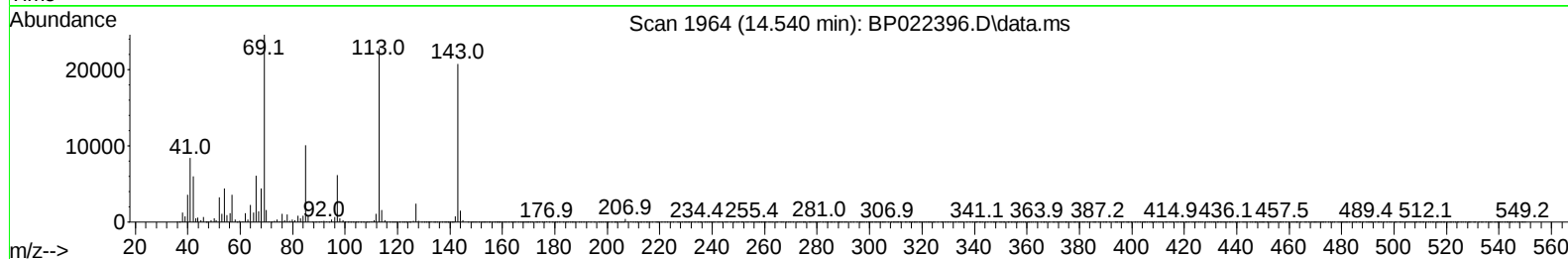
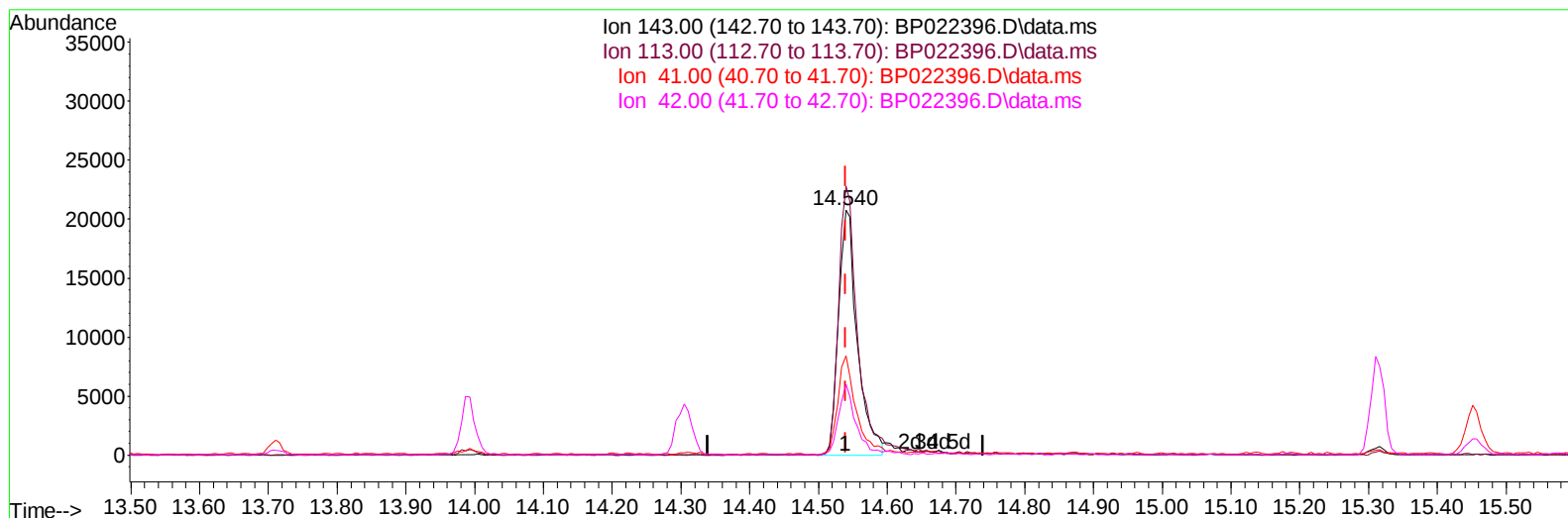
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022396.D
 Acq On : 15 Oct 2024 19:40
 Operator : RC/JU
 Sample : P4349-14
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAH0

Manual IntegrationsAPPROVED

Quant Time: Oct 16 02:34:33 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: BP022396.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (+ 0.000) 11.83 ng/u1

response 39294

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.70
41.00	35.90	40.48
42.00	23.70	28.96#

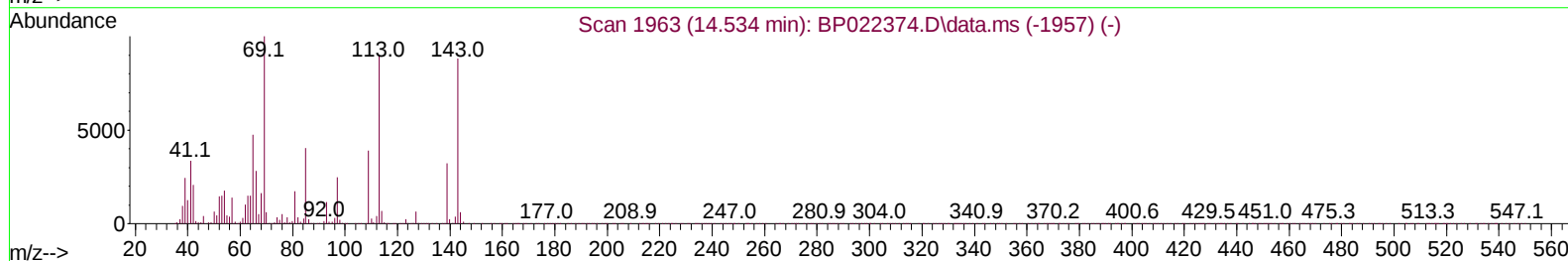
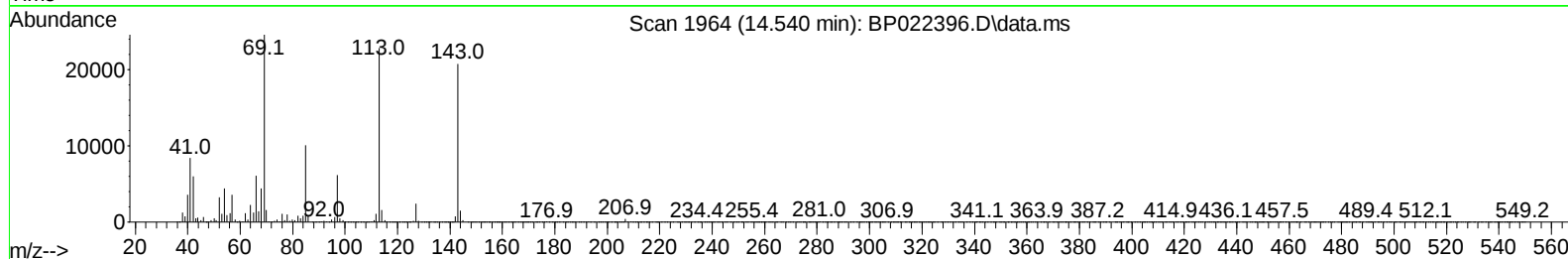
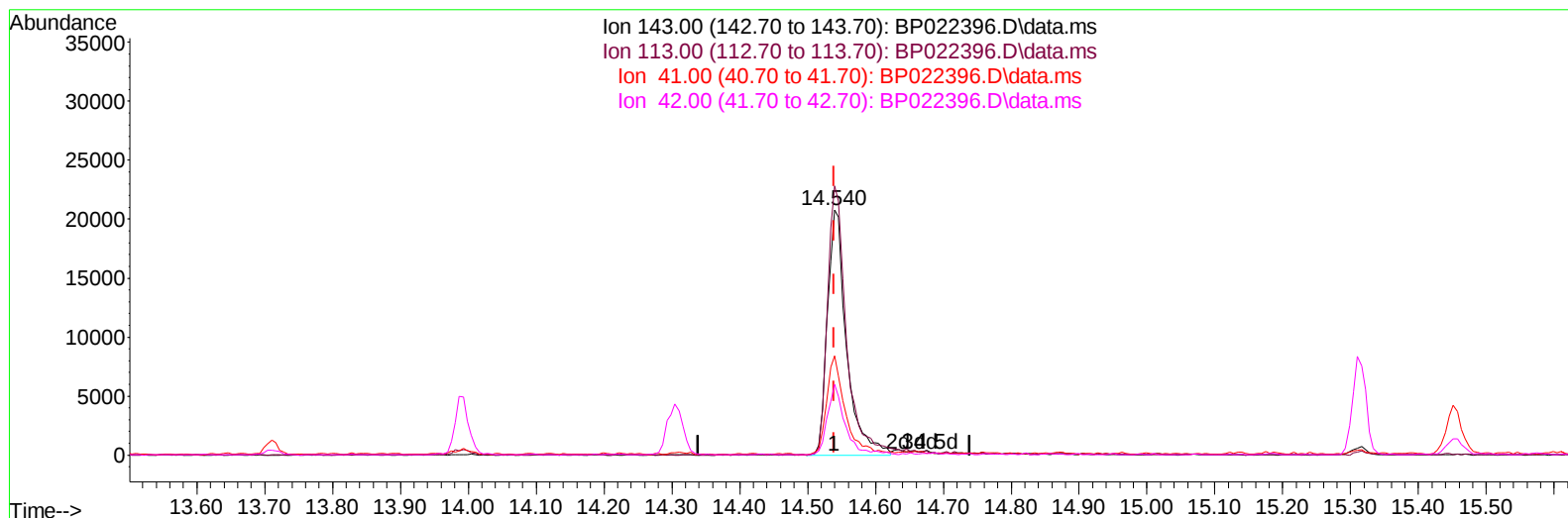
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TIC: BP022396.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (+ 0.000) 12.21 ng/ul m

response 40539

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.70
41.00	35.90	40.48
42.00	23.70	28.96#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/16/2024
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Quant Time: Oct 16 02:36:01 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	80831	20.000	ng/ul	0.00
20) Naphthalene-d8	10.452	136	311820	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.304	164	249109	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	619621	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	718234	20.000	ng/ul	# 0.00
88) Perylene-d12	24.833	264	861691	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	7963	3.828	ng/uL	0.00
4) Pyridine-d5	3.640	84	49605	8.687	ng/ul	0.00
7) Phenol-d5	6.881	99	87564	12.869	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	104714	26.185	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	134945	27.951	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	121905	22.450	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	67668	28.223	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	80022	28.829	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	166901	28.290	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	144695	20.756	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	555231	28.053	ng/ul	0.00
49) Acenaphthylene-d8	13.993	160	595587	28.332	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	40539m	12.209	ng/ul	0.00
60) Fluorene-d10	15.310	176	499200	29.079	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.457	200	106364	26.101	ng/ul	0.01
73) Anthracene-d10	17.216	188	889289	30.509	ng/ul	0.01
81) Pyrene-d10	19.586	212	1167683	30.743	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.616	264	1436896	31.224	ng/ul	-0.02

Target Compounds Qvalue

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

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