

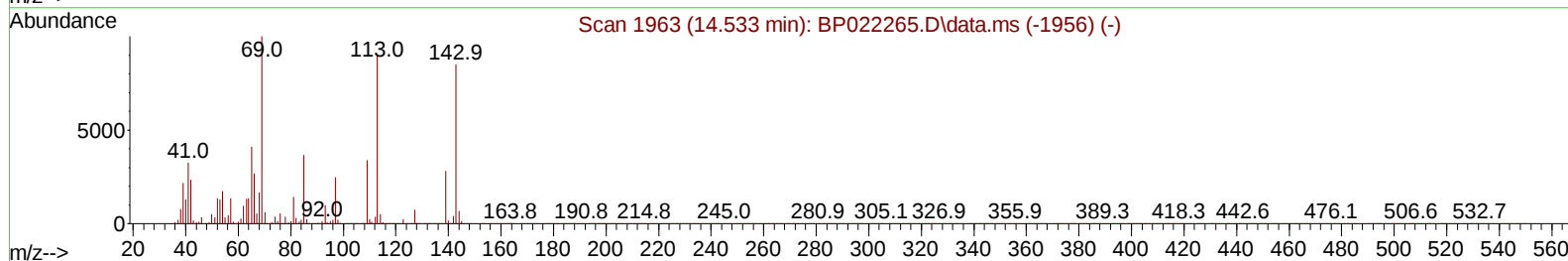
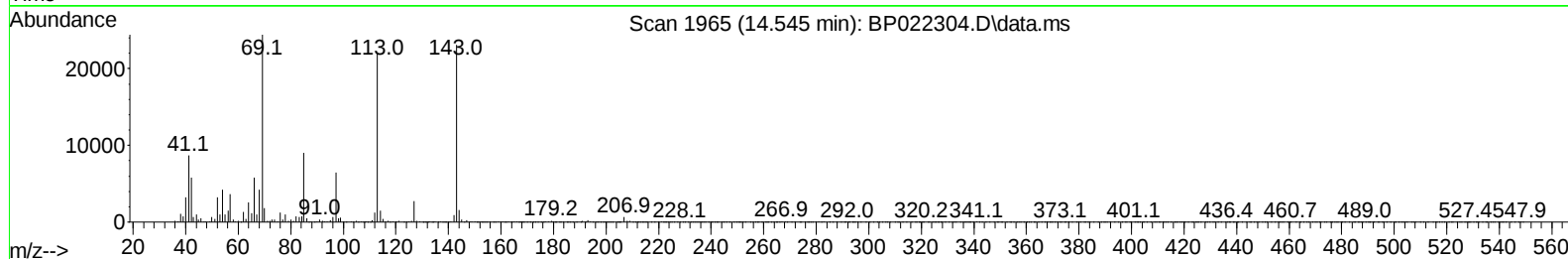
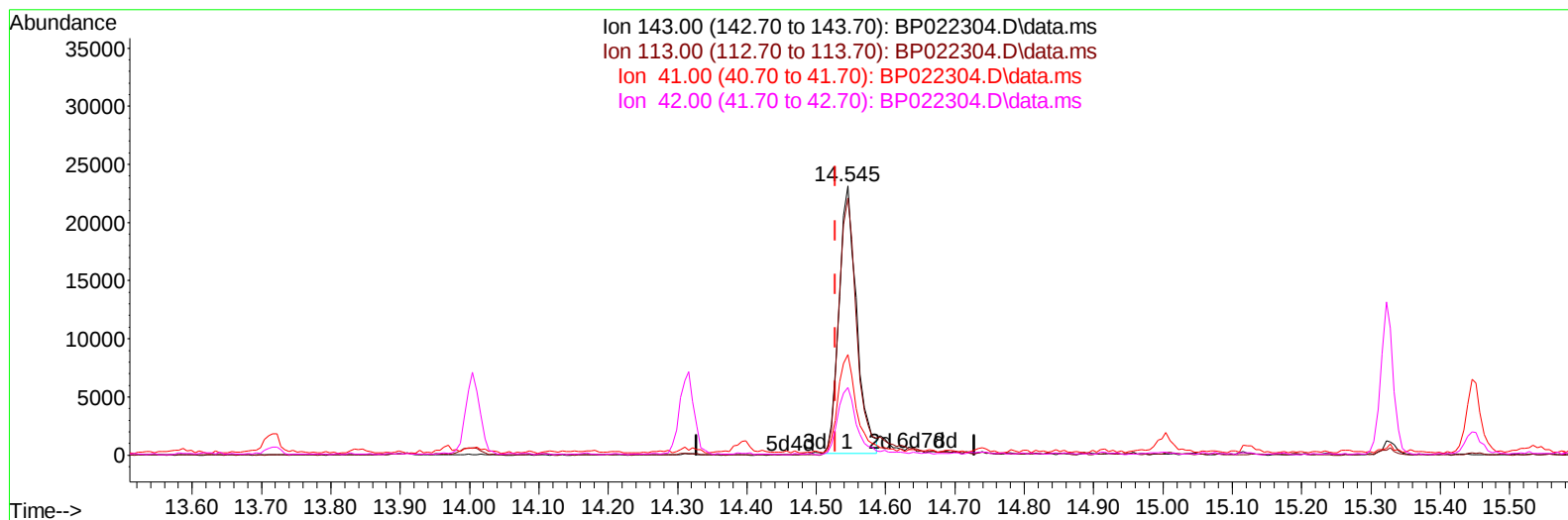
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022304.D
Acq On : 10 Oct 2024 01:13
Operator : RC/JU
Sample : P4207-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY084

Manual IntegrationsAPPROVED

Quant Time: Oct 10 02:04:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 08 02:14:47 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/10/2024
Supervised By :mohammad ahmed 10/11/2024



TIC: BP022304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.018) 7.48 ng/u1

response 40240

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	95.50
41.00	35.90	37.43
42.00	23.70	25.22

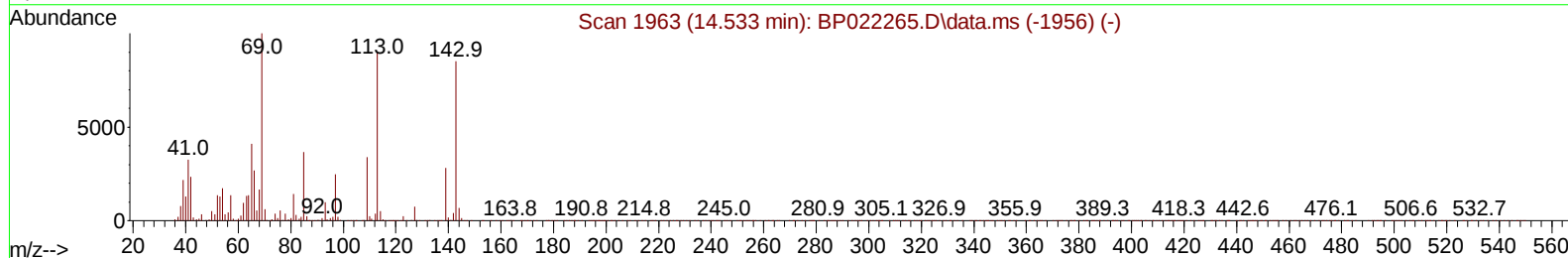
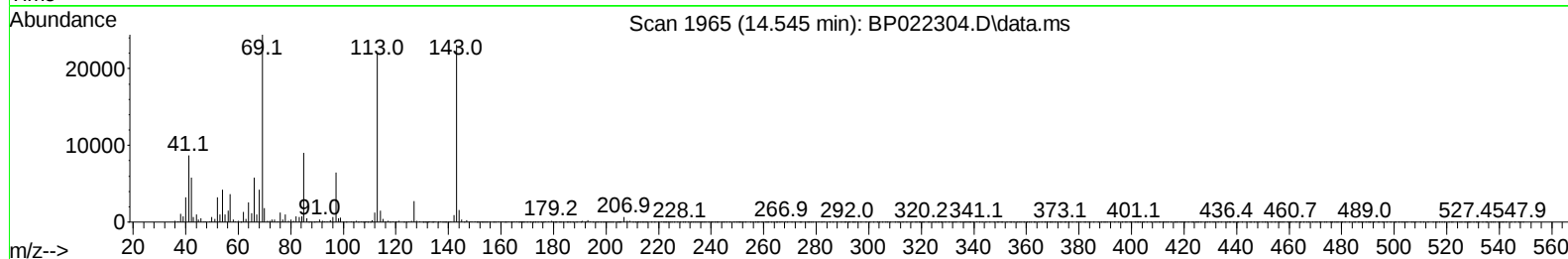
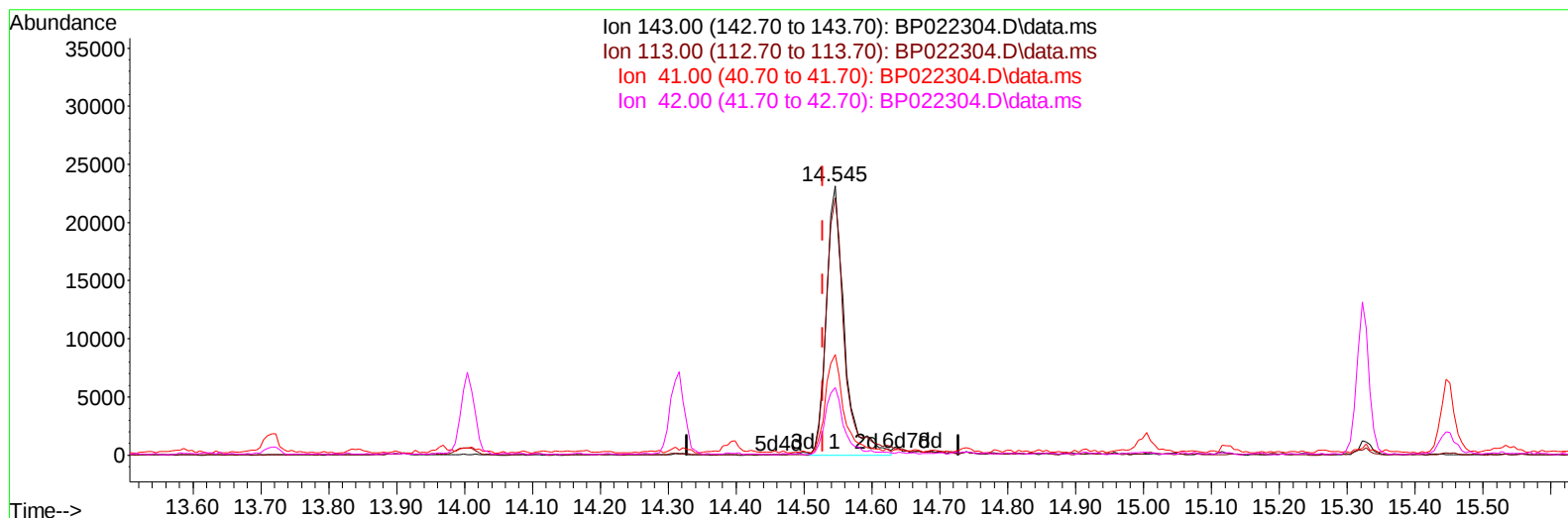
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TIC: BP022304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.018) 8.02 ng/ul m

response 43164

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	95.50
41.00	35.90	37.43
42.00	23.70	25.22

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Instrument :
 BNA_P
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 EY084

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/10/2024
 Supervised By :mohammad ahmed 10/11/2024

Quant Time: Oct 10 02:13:29 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 02:14:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	125973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	505785	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	403534	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	979510	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	963357	20.000	ng/ul	0.01
88) Perylene-d12	24.827	264	1106989	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	11835	3.651	ng/uL	0.00
4) Pyridine-d5	3.646	84	62605	7.035	ng/ul	0.01
7) Phenol-d5	6.899	99	94519	8.913	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.046	67	177109	28.417	ng/ul	0.00
11) 2-Chlorophenol-d4	7.252	132	208862	27.758	ng/ul	0.00
15) 4-Methylphenol-d8	8.410	113	171519	20.268	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	116690	30.005	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	143655	31.906	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	281879	29.456	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	300781	26.600	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	954156	29.760	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1015988	29.835	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	43164m	8.025	ng/ul	0.02
60) Fluorene-d10	15.322	176	847585	30.479	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	148059	22.983	ng/ul	0.00
73) Anthracene-d10	17.210	188	1473361	31.975	ng/ul	0.00
81) Pyrene-d10	19.592	212	1797584	35.285	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.610	264	1919972	32.476	ng/ul	0.01

Target Compounds Qvalue

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

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