

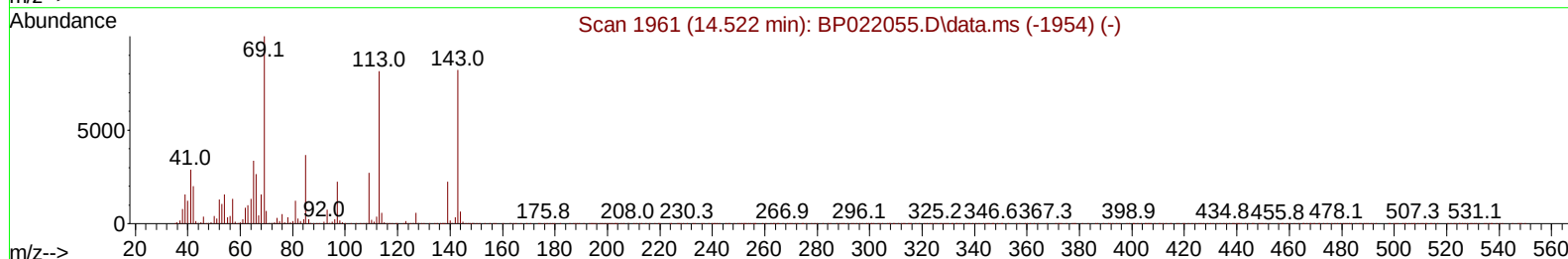
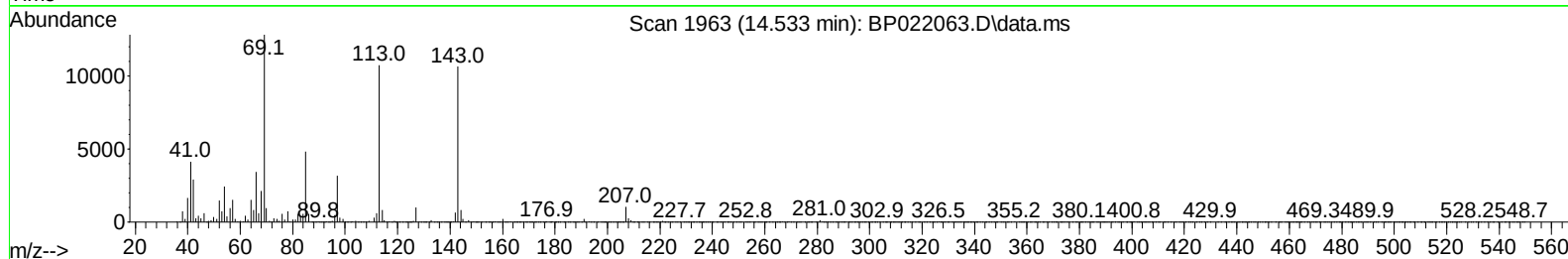
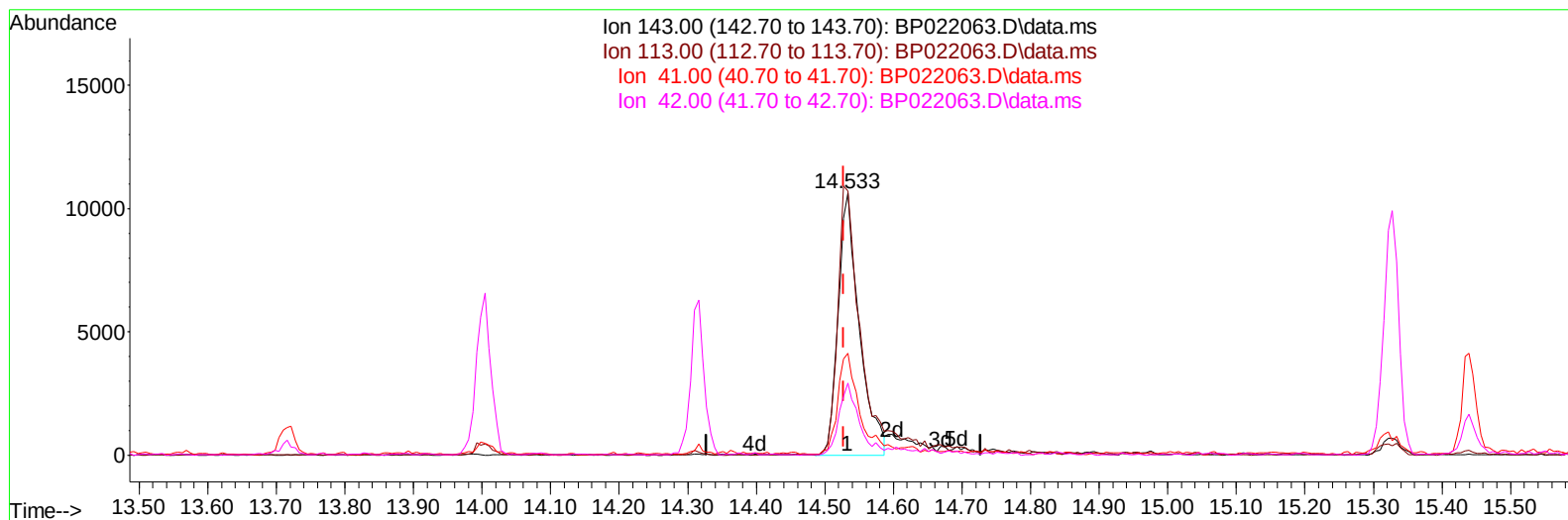
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022063.D
 Acq On : 26 Sep 2024 15:10
 Operator : RC/JU
 Sample : P4080-05
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AX4

Manual IntegrationsAPPROVED

Quant Time: Sep 27 01:17:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2024
 Supervised By :mohammad ahmed 09/28/2024



TIC: BP022063.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.533min (+ 0.006) 7.08 ng/ul

response 22605

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	100.77
41.00	35.80	38.85
42.00	26.20	27.59

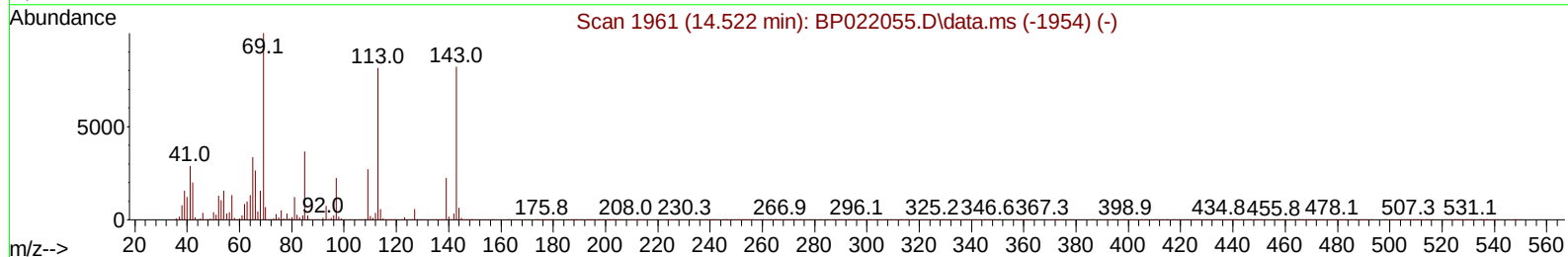
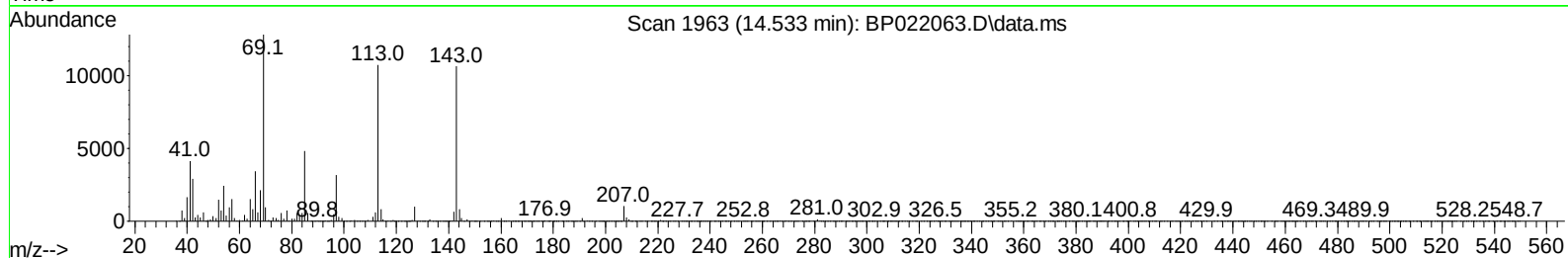
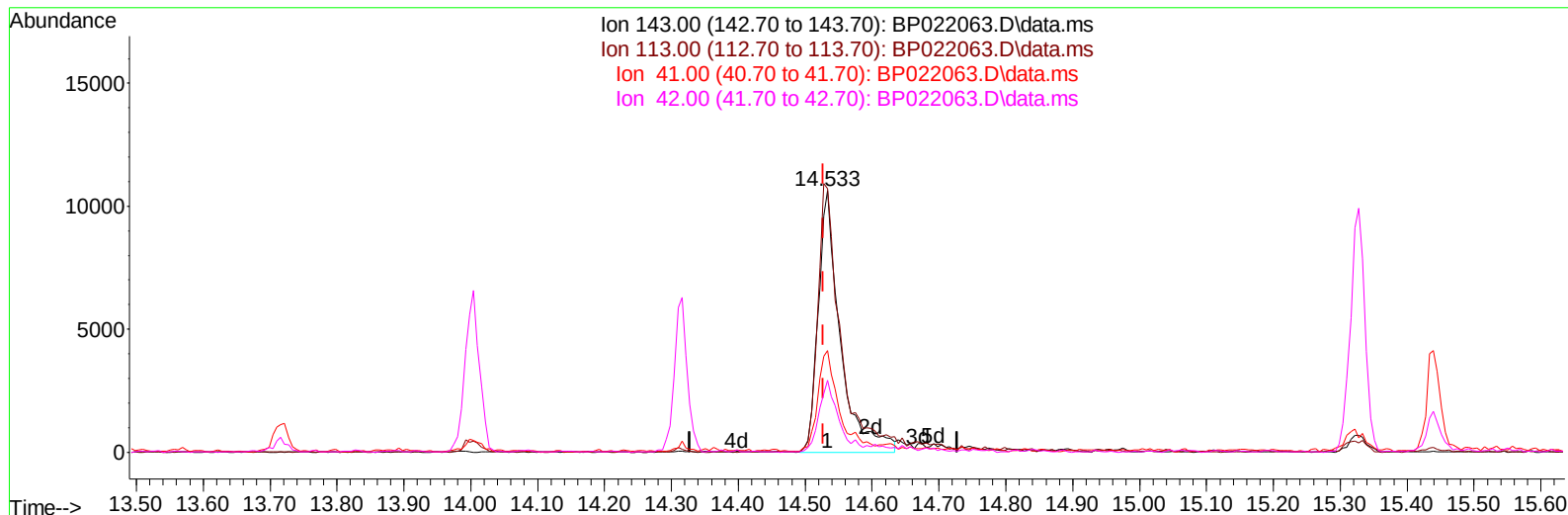
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(54) 4-Nitrophenol-d4 (S)

14.533min (+ 0.006) 7.65 ng/ul m

response 24442

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	100.77
41.00	35.80	38.85
42.00	26.20	27.59

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	88778	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	327435	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	241489	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	593192	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	685510	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264	841853	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	10760	4.747	ng/uL	0.00
4) Pyridine-d5	3.634	84	55171	8.516	ng/ul	0.00
7) Phenol-d5	6.893	99	66457	9.151	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	148729	33.260	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	166618	31.711	ng/ul	0.00
15) 4-Methylphenol-d8	8.398	113	125094	21.738	ng/ul	0.00
21) Nitrobenzene-d5	8.845	128	87923	35.876	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	102803	36.996	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.098	165	207015	34.305	ng/ul	0.00
31) 4-Chloroaniline-d4	10.598	131	193986	26.942	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	689444	36.933	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	702894	35.564	ng/ul	0.00
54) 4-Nitrophenol-d4	14.533	143	24442m	7.655	ng/ul	0.00
60) Fluorene-d10	15.327	176	600299	36.354	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	99262	26.953	ng/ul	0.00
73) Anthracene-d10	17.210	188	1085090	39.184	ng/ul	0.00
81) Pyrene-d10	19.598	212	1435232	40.412	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.603	264	1755872	39.655	ng/ul	0.00

Target Compounds Qvalue

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

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