

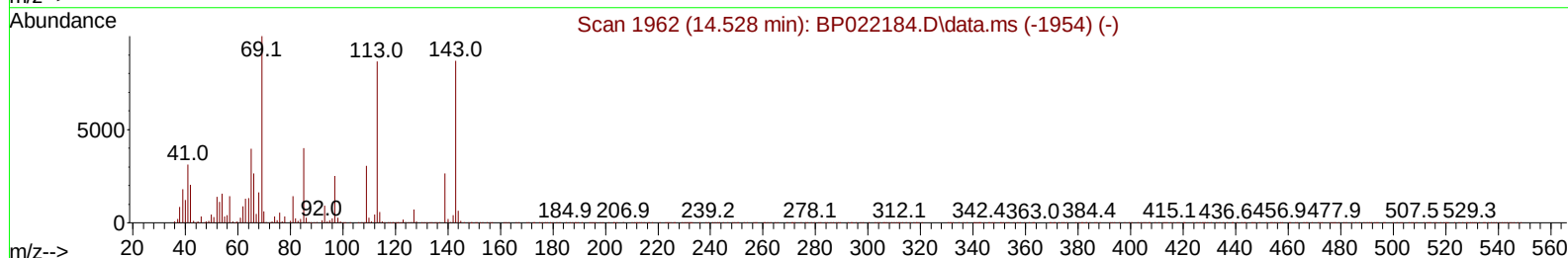
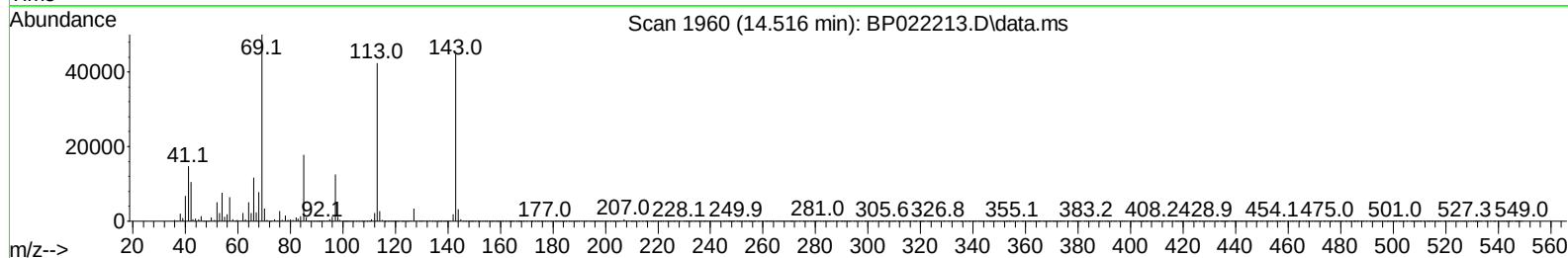
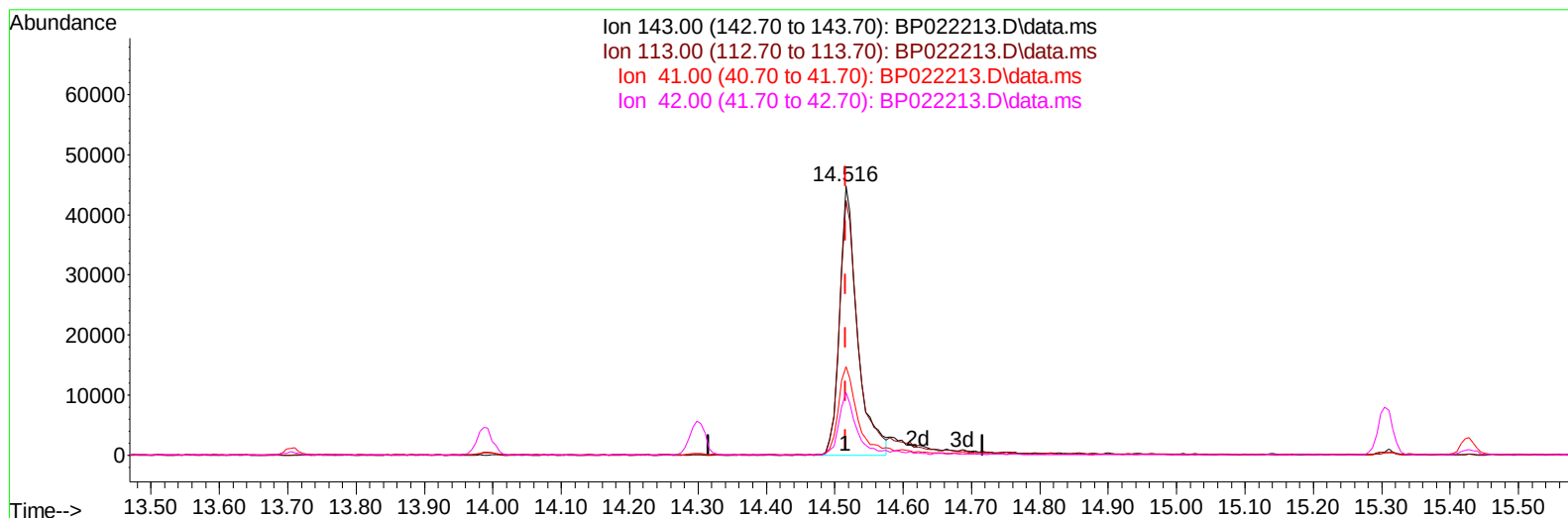
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022213.D
Acq On : 04 Oct 2024 04:01
Operator : RC/JU
Sample : PB163860BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK860

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
Supervised By :mohammad ahmed 10/05/2024

Quant Time: Oct 04 05:28:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 04 05:25:28 2024
Response via : Initial Calibration



TIC: BP022213.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.000) 17.89 ng/ul

response 81099

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	94.57
41.00	35.80	32.93
42.00	26.20	23.35

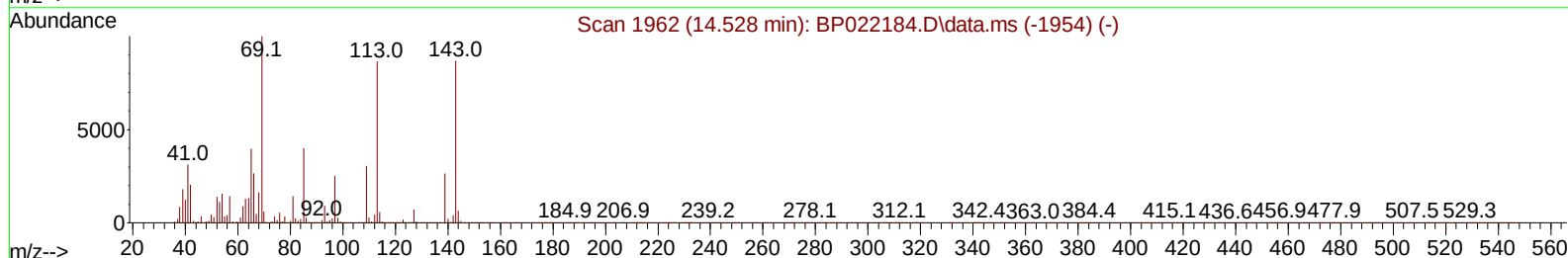
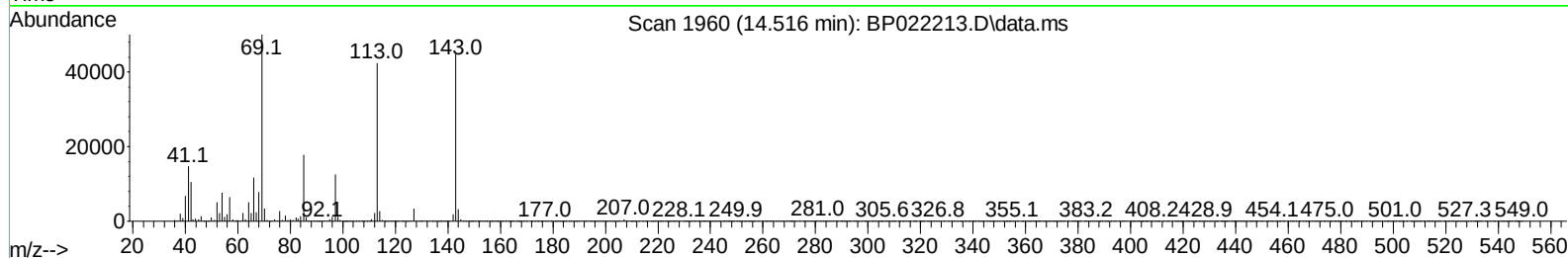
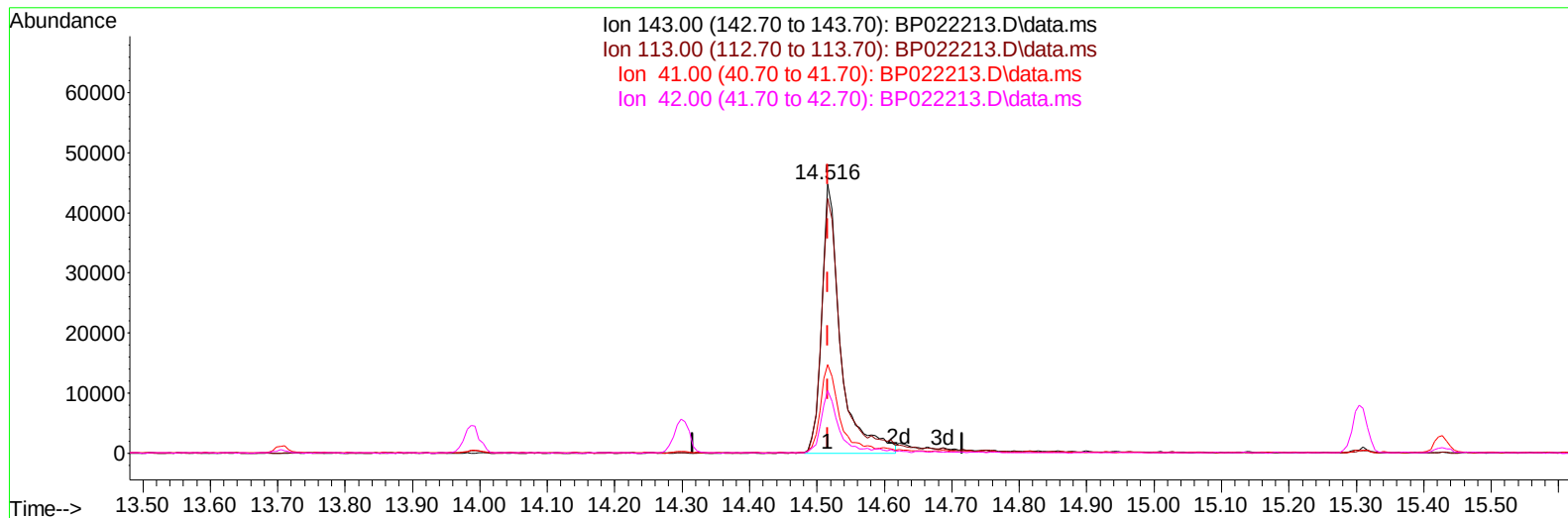
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TIC: BP022213.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.000) 19.09 ng/ul m

response 86561

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	94.57
41.00	35.80	32.93
42.00	26.20	23.35

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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 05:25:28 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	112843	20.000	ng/ul	0.00
20) Naphthalene-d8	10.452	136	437821	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164	342869	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	827340	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	877620	20.000	ng/ul	-0.01
88) Perylene-d12	24.798	264	976761	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	12564	4.361	ng/uL	0.00
4) Pyridine-d5	3.628	84	193147	23.455	ng/ul	0.00
7) Phenol-d5	6.881	99	220311	23.867	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	131196	23.082	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	175797	26.323	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	178480	24.401	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	83417	25.456	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	100341	27.006	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	202892	25.145	ng/ul	0.00
31) 4-Chloroaniline-d4	10.581	131	239597	24.887	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	679935	25.654	ng/ul	0.00
49) Acenaphthylene-d8	13.987	160	733948	26.155	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	86561m	19.093	ng/ul	0.00
60) Fluorene-d10	15.310	176	599144	25.556	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.428	200	77807	15.148	ng/ul	0.00
73) Anthracene-d10	17.192	188	997778	25.833	ng/ul	-0.01
81) Pyrene-d10	19.575	212	1228876	27.027	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.574	264	1331666	25.921	ng/ul	0.00

Target Compounds Qvalue

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

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