

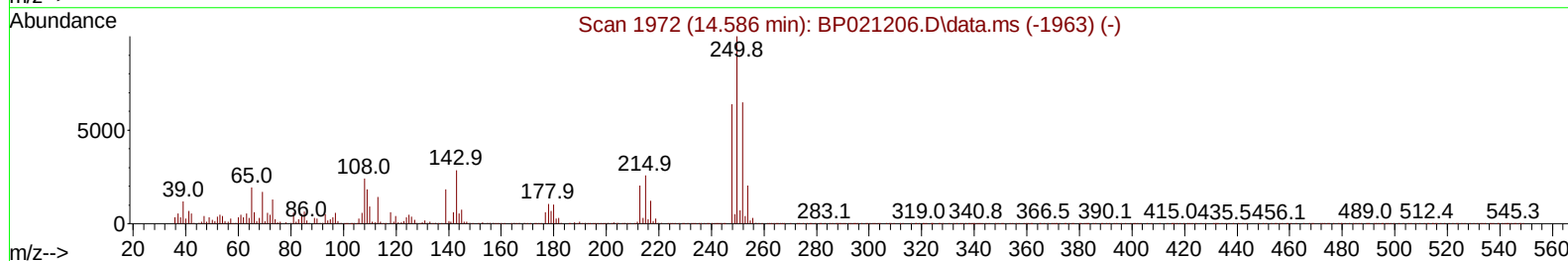
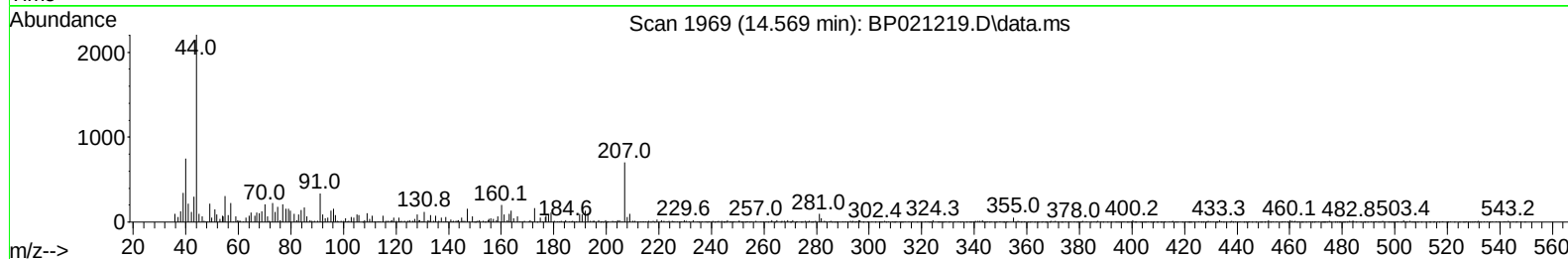
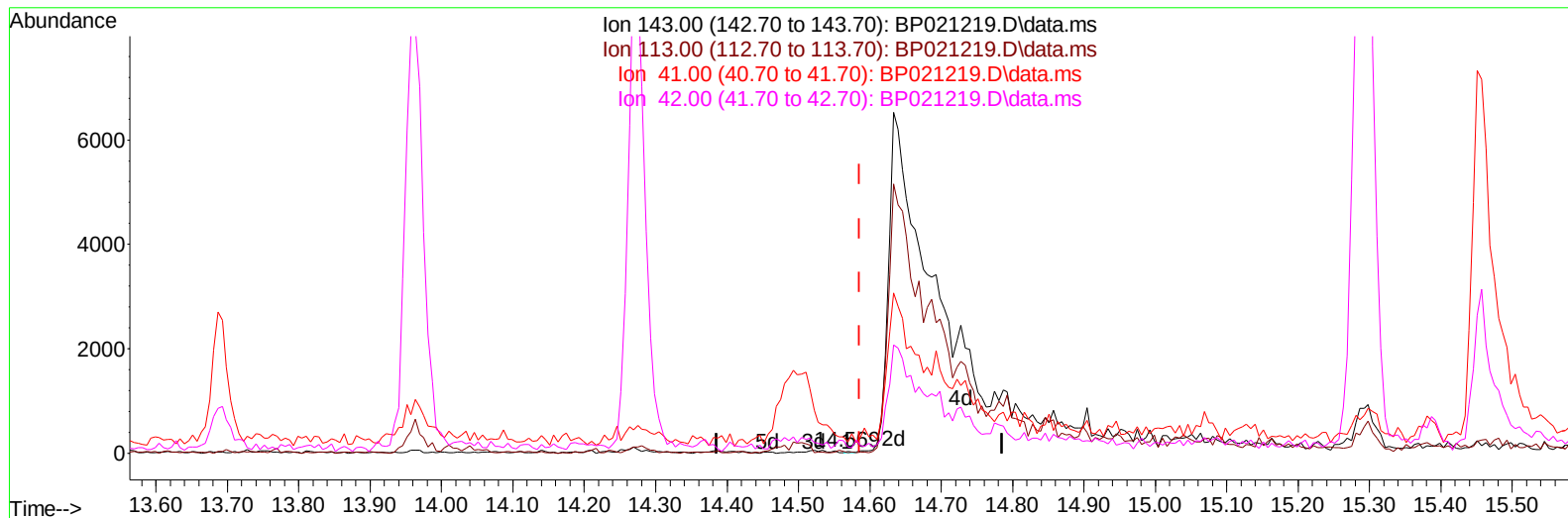
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021219.D
Acq On : 29 Jul 2024 22:13
Operator : MA/JU
Sample : P3280-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ1

Manual IntegrationsAPPROVED

Quant Time: Jul 29 22:48:20 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/30/2024
Supervised By :mohammad ahmed 09/04/2024



TIC: BP021219.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.017) 0.00 ng/u1

response 19

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	42.31#
41.00	37.90	846.15#
42.00	27.10	465.38#

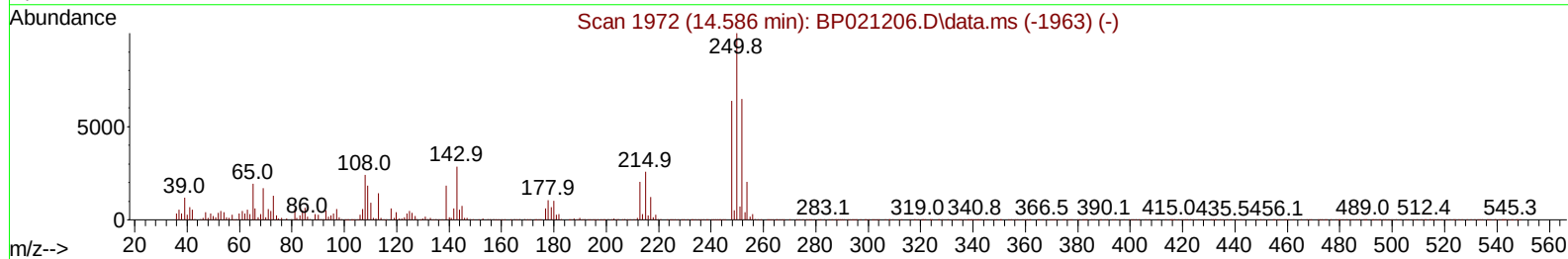
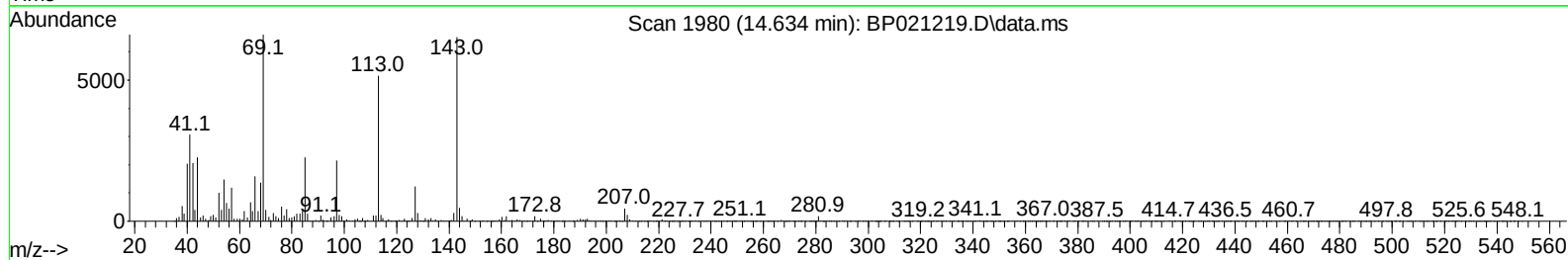
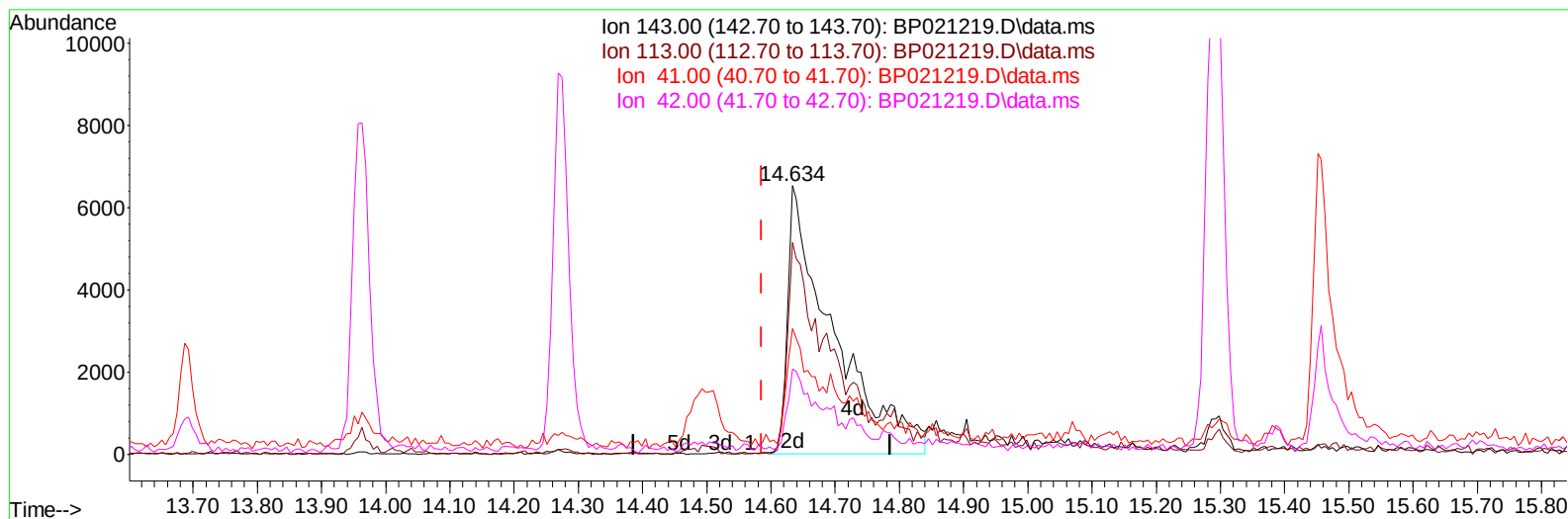
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TIC: BP021219.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.048) 5.99 ng/ul m

response 31955

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	78.86
41.00	37.90	46.91#
42.00	27.10	31.70

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/30/2024
 Supervised By :mohammad ahmed 09/04/2024

Quant Time: Jul 29 23:01:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	201631	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	801955	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	515399	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1111918	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1114446	20.000	ng/ul	0.00
88) Perylene-d12	24.839	264	1343568	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	20798	4.694	ng/uL	0.00
4) Pyridine-d5	3.605	84	101782	7.935	ng/ul	0.01
7) Phenol-d5	6.869	99	104783	6.374	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67	276875	27.793	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	325529	24.035	ng/ul	0.01
15) 4-Methylphenol-d8	8.375	113	216303	15.843	ng/ul	0.01
21) Nitrobenzene-d5	8.799	128	173797	27.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.511	143	199537	27.988	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.058	165	333178	25.845	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	371617	21.760	ng/ul	0.01
46) Dimethylphthalate-d6	13.693	166	1160124	30.964	ng/ul	0.01
49) Acenaphthylene-d8	13.963	160	1241714	29.515	ng/ul	0.00
54) 4-Nitrophenol-d4	14.634	143	31955m	5.994	ng/ul	0.05
60) Fluorene-d10	15.293	176	977566	31.804	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	185719	27.605	ng/ul	0.00
73) Anthracene-d10	17.198	188	1704654	34.517	ng/ul	0.00
81) Pyrene-d10	19.581	212	2123806	30.825	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	2453822	37.031	ng/ul	0.00

Target Compounds Qvalue

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

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