

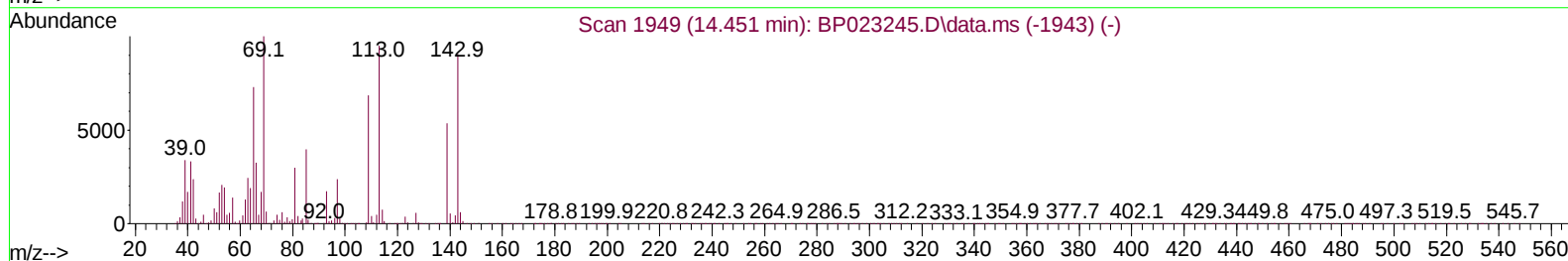
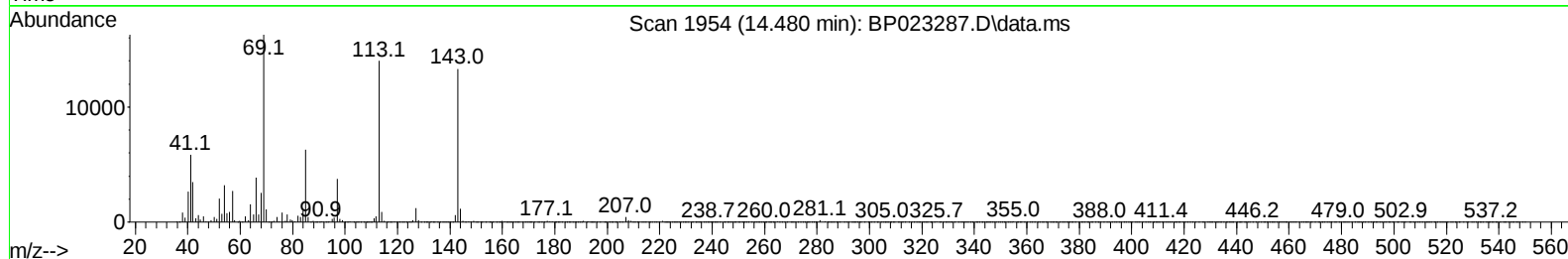
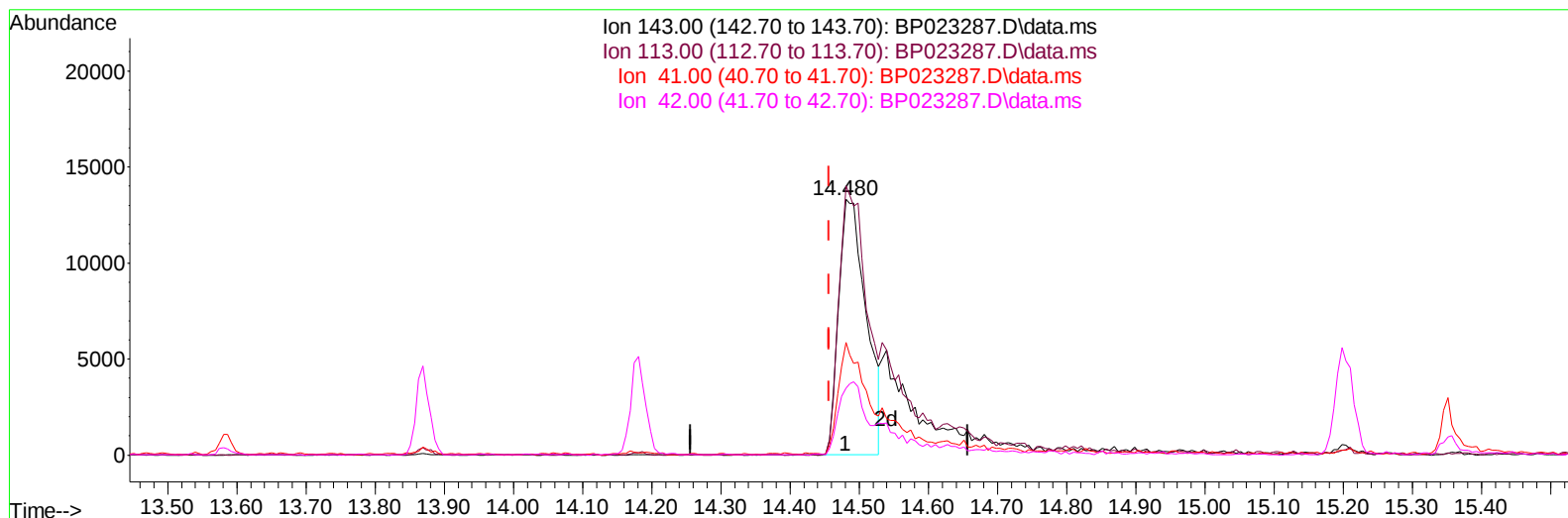
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023287.D
Acq On : 21 Nov 2024 16:44
Operator : RC/JU
Sample : PB165140BL
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK140

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/22/2024
Supervised By :mohammad ahmed 11/22/2024

Quant Time: Nov 22 00:45:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 20 04:34:35 2024
Response via : Initial Calibration



TIC: BP023287.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.480min (+ 0.023) 13.63 ng/u1****response 36568**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.14
41.00	39.20	43.97
42.00	25.90	26.13

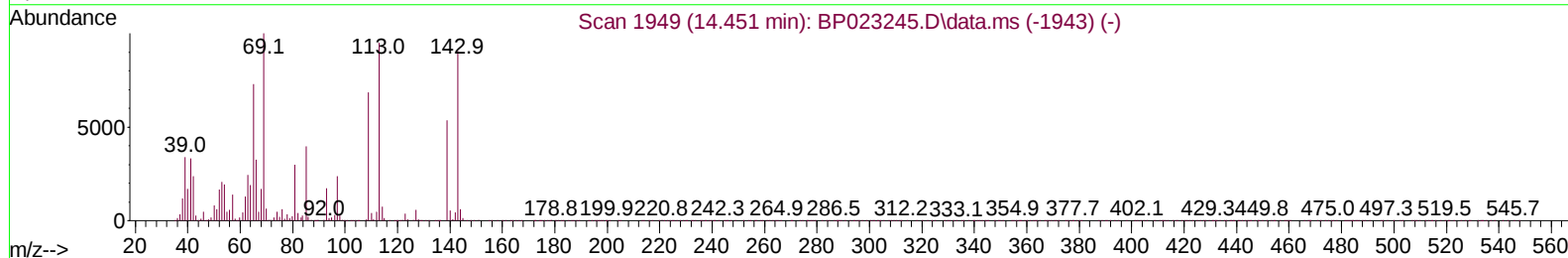
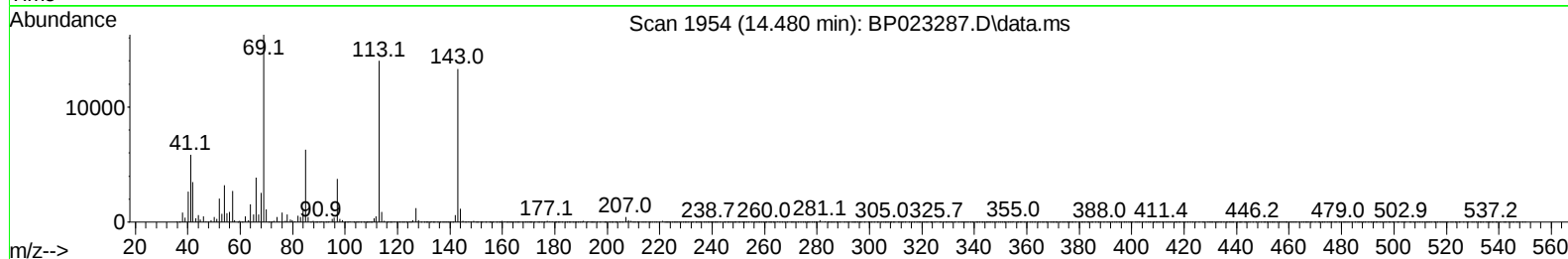
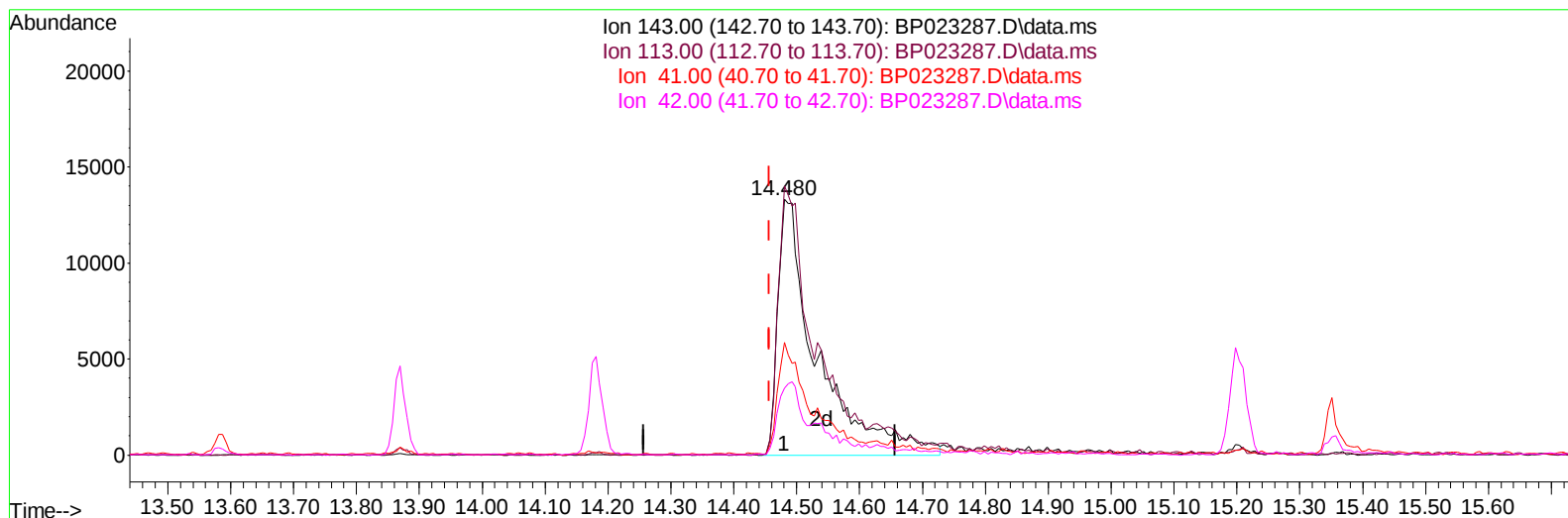
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TIC: BP023287.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.480min (+ 0.023) 21.51 ng/ul m

response 57695

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.14
41.00	39.20	43.97
42.00	25.90	26.13

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	86151	20.000	ng/ul	0.00
20) Naphthalene-d8	10.310	136	323684	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.181	164	247659	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	596611	20.000	ng/ul	0.00
79) Chrysene-d12	21.439	240	623665	20.000	ng/ul	0.02
88) Perylene-d12	24.668	264	697679	20.000	ng/ul	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	9941	5.547	ng/uL	0.00
4) Pyridine-d5	3.552	84	123758	23.241	ng/ul	0.00
7) Phenol-d5	6.769	99	156519	24.574	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.910	67	97666	26.099	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	119820	25.686	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	129376	24.683	ng/ul	0.00
21) Nitrobenzene-d5	8.704	128	56783	25.199	ng/ul	0.00
24) 2-Nitrophenol-d4	9.416	143	69285	25.790	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	136300	24.227	ng/ul	0.00
31) 4-Chloroaniline-d4	10.463	131	152080	22.659	ng/ul	0.00
46) Dimethylphthalate-d6	13.581	166	480068	26.895	ng/ul	0.00
49) Acenaphthylene-d8	13.869	160	499013	26.793	ng/ul	0.00
54) 4-Nitrophenol-d4	14.480	143	57695m	21.508	ng/ul	0.02
60) Fluorene-d10	15.204	176	405989	25.943	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.351	200	63248	17.697	ng/ul	0.00
73) Anthracene-d10	17.092	188	686850	27.190	ng/ul	0.00
81) Pyrene-d10	19.486	212	862854	30.037	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	932788	28.068	ng/ul	0.02

Target Compounds Qvalue

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

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