

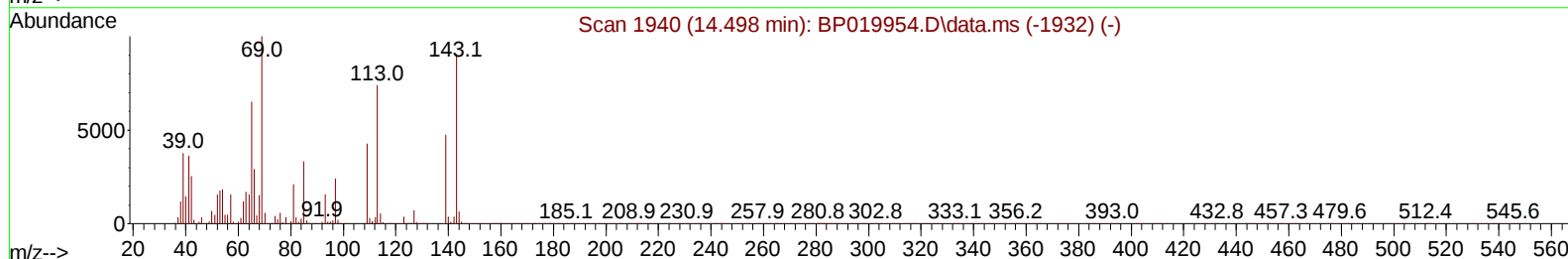
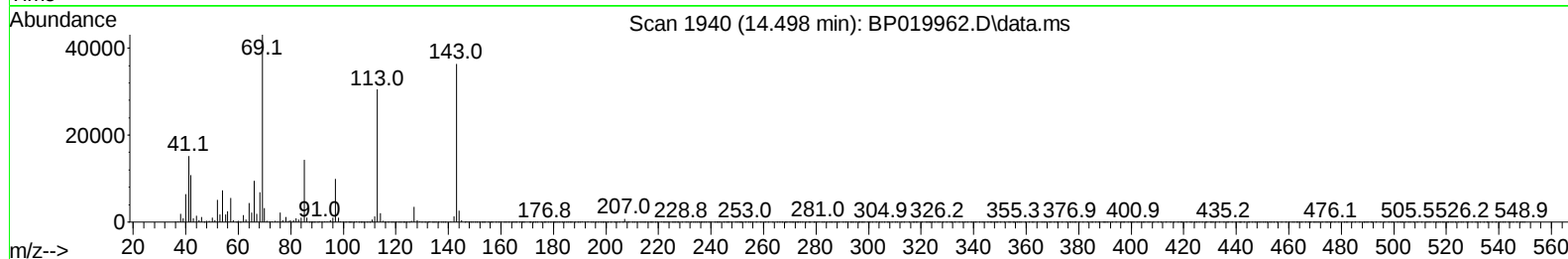
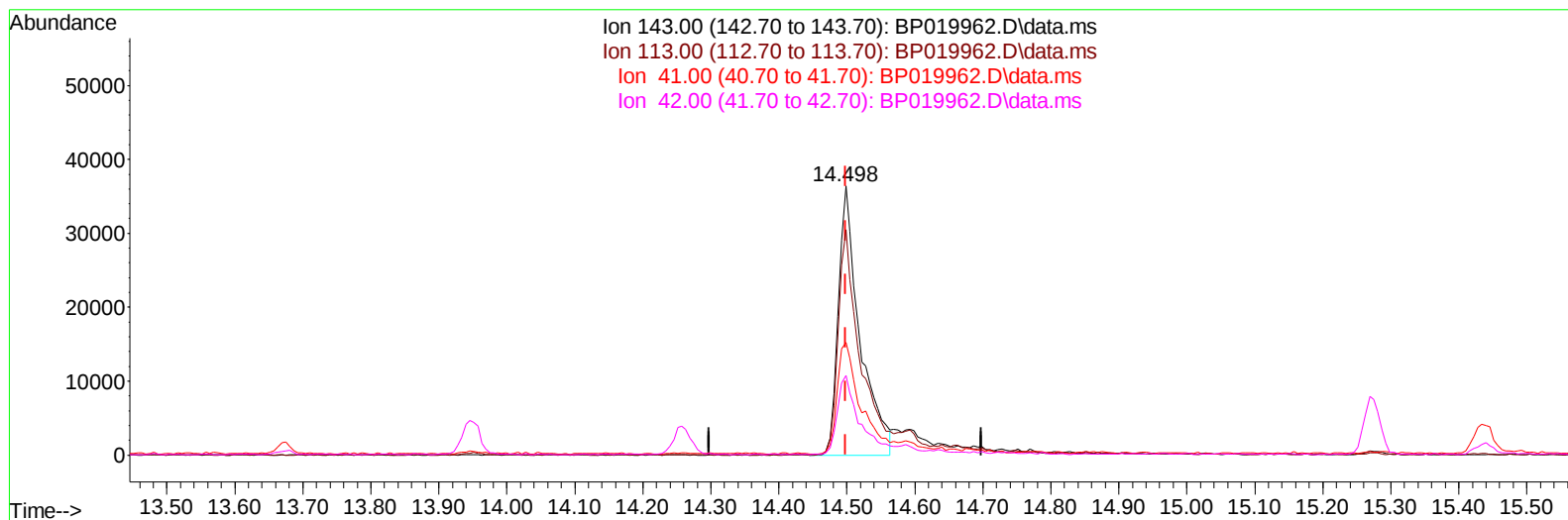
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019962.D
Acq On : 22 Mar 2024 19:15
Operator : MA/JU
Sample : PB159746BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK746

Manual IntegrationsAPPROVED

Quant Time: Mar 22 23:03:10 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 22 22:40:53 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/26/2024
Supervised By :mohammad ahmed 03/26/2024



TIC: BP019962.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.000) 31.43 ng/u1

response 79975

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	83.71
41.00	40.90	41.75
42.00	28.50	29.57

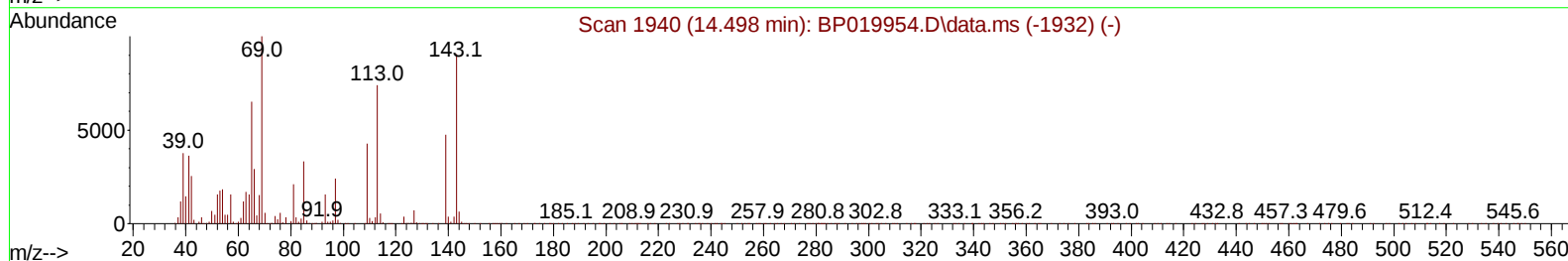
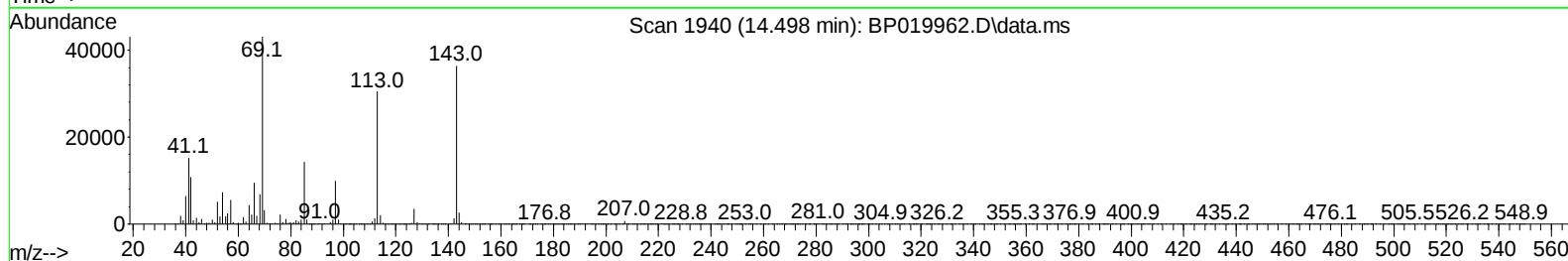
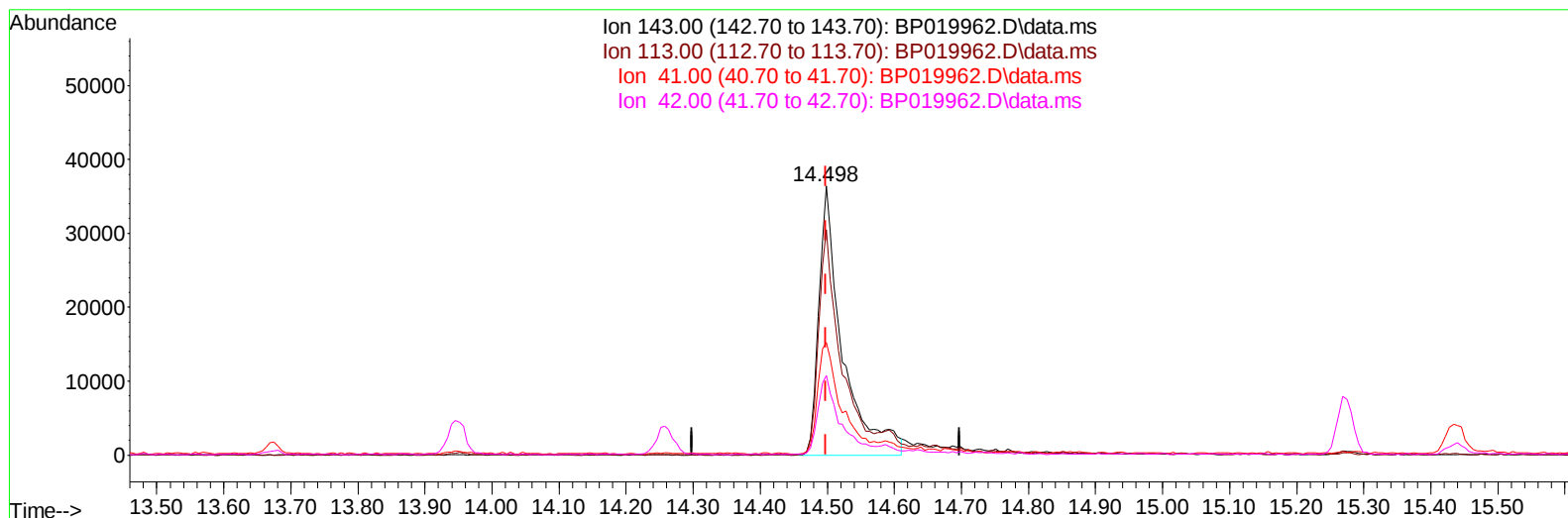
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TIC: BP019962.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.000) 34.86 ng/ul m

response 88713

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	83.71
41.00	40.90	41.75
42.00	28.50	29.57

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/26/2024
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Quant Time: Mar 22 23:19:35 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 22 22:40:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	91015	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.375	136	353856	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	230888	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	499695	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	500815	20.000	ng/ul	0.00
88) Perylene-d12	24.863	264	569655	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	16565	7.974	ng/uL	0.00
4) Pyridine-d5	3.640	84	207433	33.071	ng/ul	0.00
7) Phenol-d5	6.799	99	272126	35.451	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.975	67	178583	35.709	ng/ul	0.00
11) 2-Chlorophenol-d4	7.152	132	198304	35.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	208012	35.026	ng/ul	-0.01
21) Nitrobenzene-d5	8.763	128	98092	36.167	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.469	143	111359	36.696	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	187060	33.753	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	246817	34.162	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	637274	39.367	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	688828	36.426	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	88713m	34.859	ng/ul	0.00
60) Fluorene-d10	15.275	176	520444	37.741	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.440	200	98170	32.821	ng/ul	0.00
73) Anthracene-d10	17.187	188	870323	39.346	ng/ul	0.00
81) Pyrene-d10	19.581	212	1076345	36.631	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.616	264	1111486	40.614	ng/ul	-0.02

Target Compounds Qvalue

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

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