

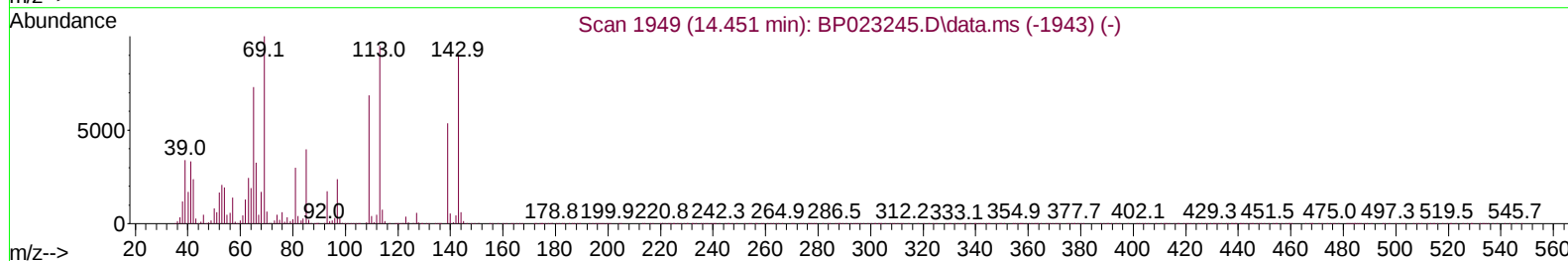
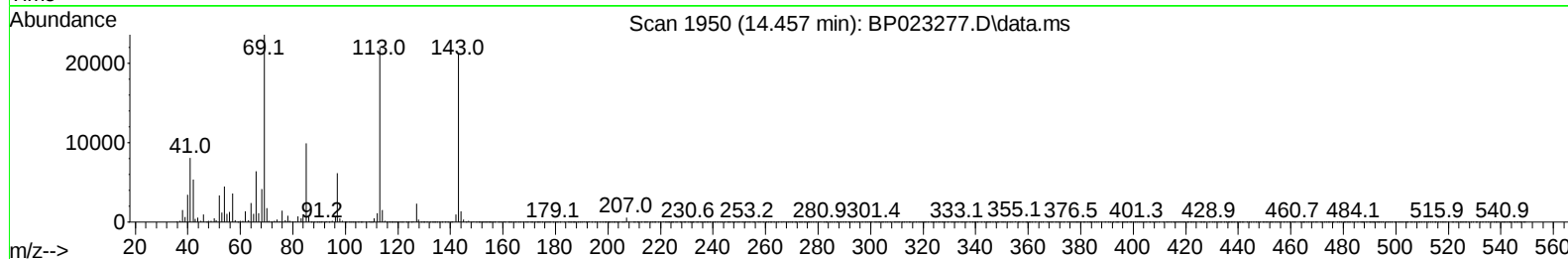
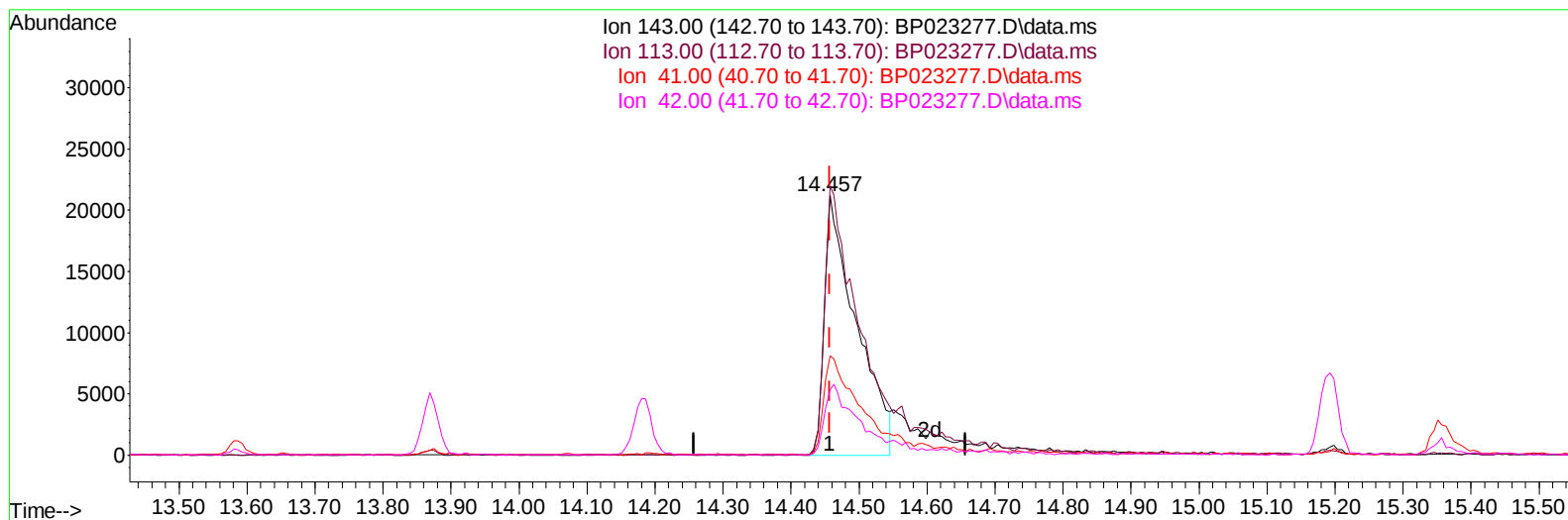
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023277.D
Acq On : 21 Nov 2024 09:57
Operator : RC/JU
Sample : PB165090BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK090

Manual IntegrationsAPPROVED

Quant Time: Nov 21 10:33:05 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

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



TIC: BP023277.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.457min (+ 0.000) 23.17 ng/u1

response 68771

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	103.66
41.00	39.20	38.18
42.00	25.90	25.21

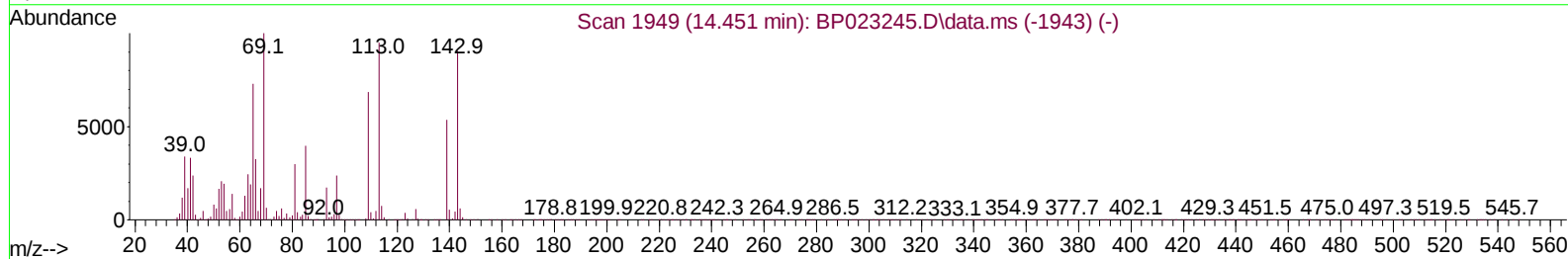
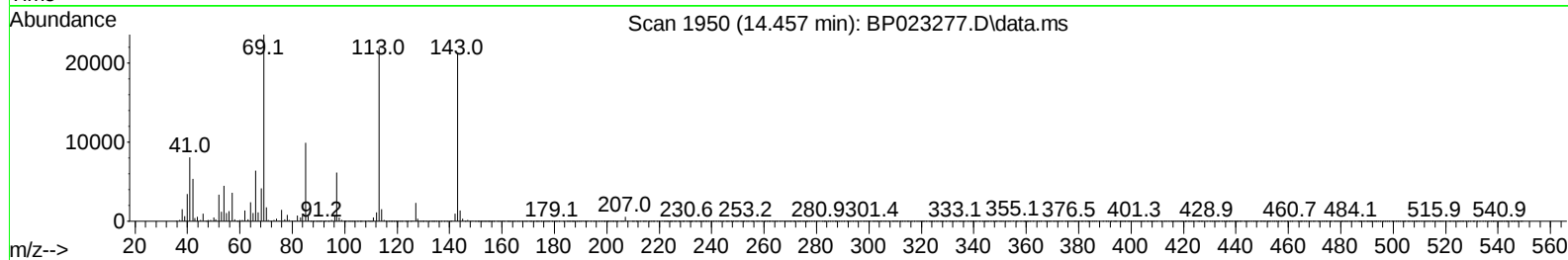
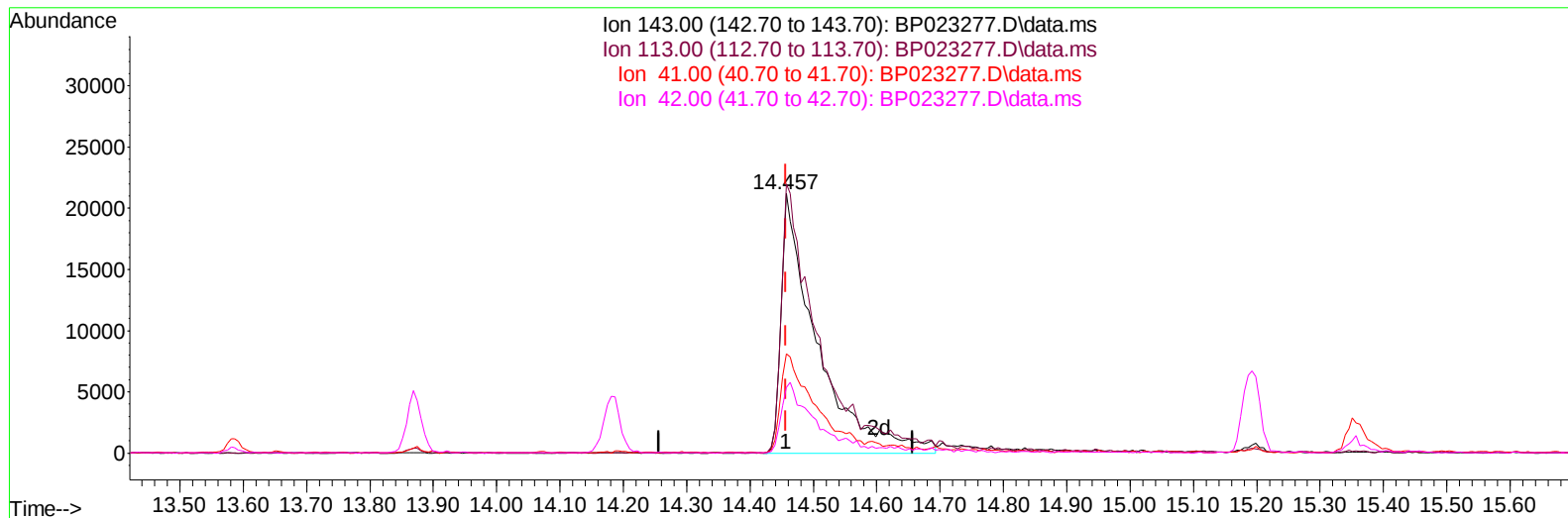
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TIC: BP023277.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.457min (+ 0.000) 28.03 ng/ul m

response 83164

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	103.66
41.00	39.20	38.18
42.00	25.90	25.21

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 BNA_P
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 SBLK090

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/22/2024
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Quant Time: Nov 22 00:04:59 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	105173	20.000	ng/ul	0.00
20) Naphthalene-d8	10.304	136	371335	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.186	164	273974	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.980	188	661392	20.000	ng/ul	-0.02
79) Chrysene-d12	21.439	240	775388	20.000	ng/ul	0.02
88) Perylene-d12	24.662	264	904714	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	14050	6.421	ng/uL	0.00
4) Pyridine-d5	3.546	84	179613	27.630	ng/ul	0.00
7) Phenol-d5	6.769	99	215631	27.731	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.916	67	133583	29.241	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	169675	29.795	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	172906	27.022	ng/ul	0.00
21) Nitrobenzene-d5	8.710	128	79886	30.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.416	143	96610	31.347	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.957	165	190323	29.488	ng/ul	0.00
31) 4-Chloroaniline-d4	10.463	131	206489	26.817	ng/ul	0.00
46) Dimethylphthalate-d6	13.586	166	622836	31.542	ng/ul	0.00
49) Acenaphthylene-d8	13.869	160	660843	32.074	ng/ul	0.00
54) 4-Nitrophenol-d4	14.457	143	83164m	28.025	ng/ul	0.00
60) Fluorene-d10	15.192	176	537165	31.028	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.351	200	92897	23.447	ng/ul	0.00
73) Anthracene-d10	17.098	188	927698	33.127	ng/ul	0.00
81) Pyrene-d10	19.480	212	1186257	33.215	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1445417	33.540	ng/ul	0.02

Target Compounds Qvalue

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

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