

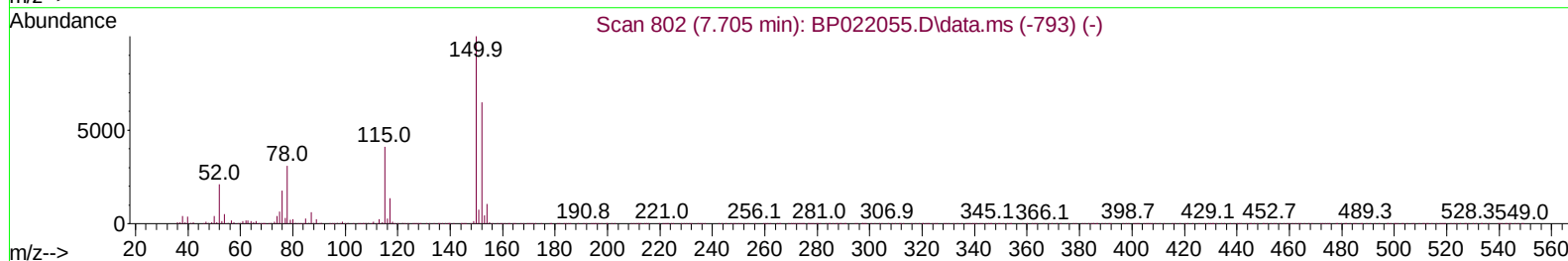
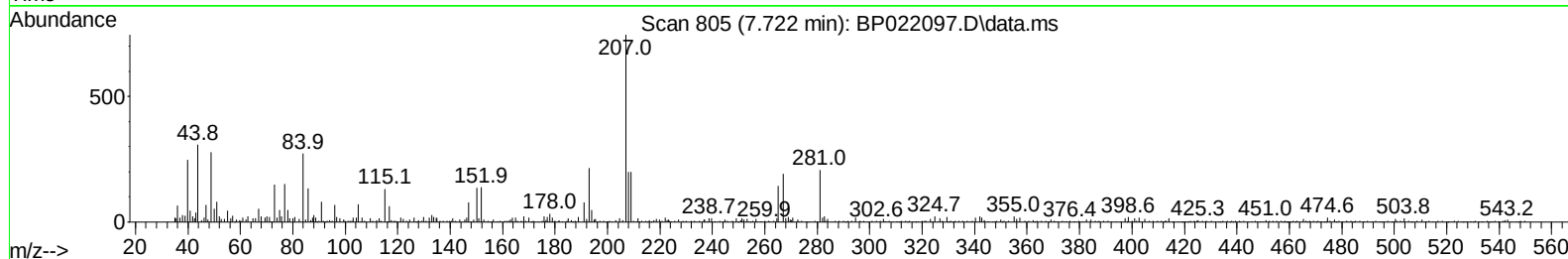
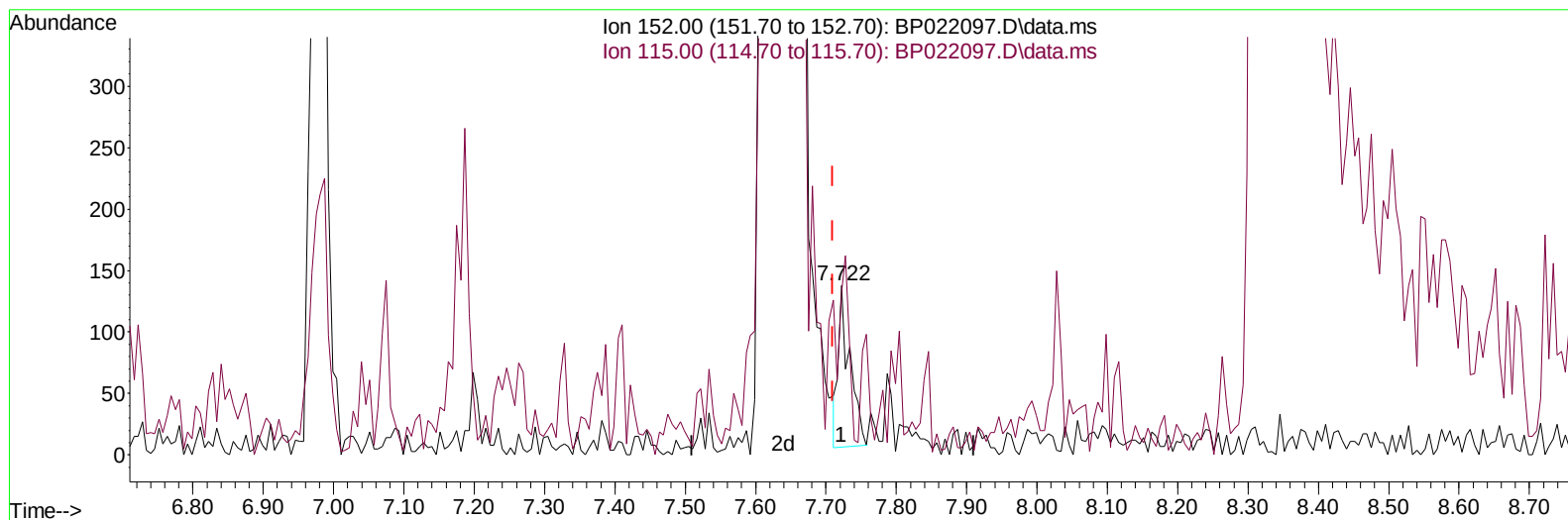
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022097.D
Acq On : 28 Sep 2024 09:38
Operator : RC/JU
Sample : PB163765BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK765

Manual IntegrationsAPPROVED

Quant Time: Sep 28 23:24:52 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/01/2024
Supervised By :mohammad ahmed 10/01/2024



TIC: BP022097.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.722min (+ 0.012) 20.00 ng/u1

response 149

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	64.00	95.65#
0.00	0.00	0.00
0.00	0.00	0.00

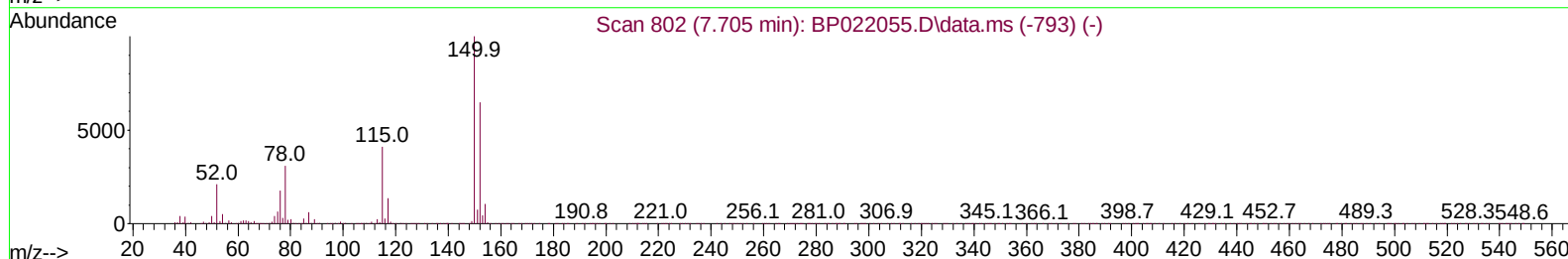
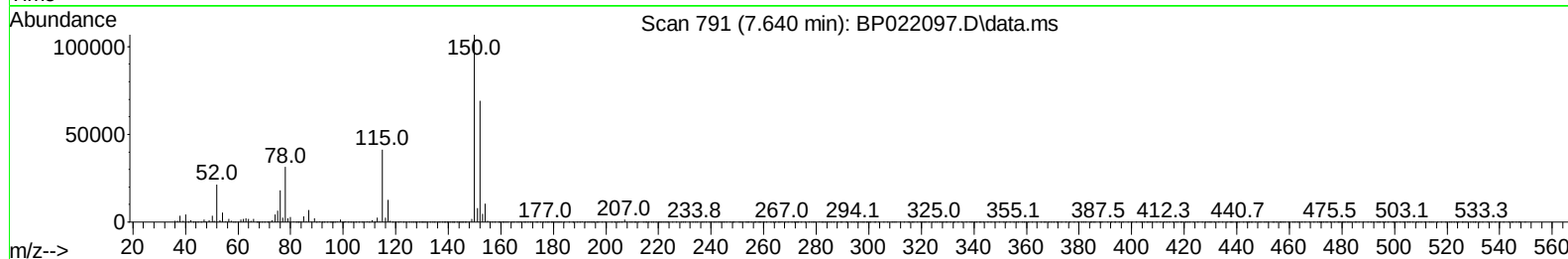
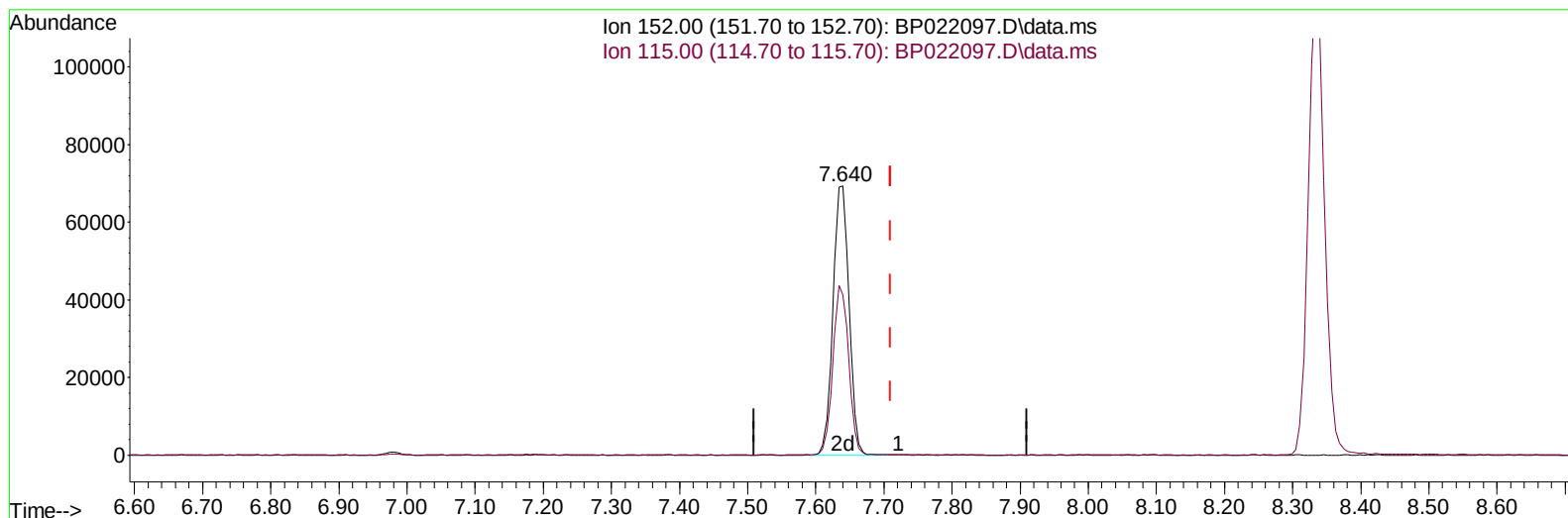
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TIC: BP022097.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.640min (-0.070) 20.00 ng/ul m

response 113096

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	64.00	59.54
0.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/01/2024
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Quant Time: Sep 28 23:26:35 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.640	152		113096m	20.000	ng/ul	-0.07
20) Naphthalene-d8	10.404	136		419487	20.000	ng/ul	-0.06
38) Acenaphthene-d10	14.286	164		316310	20.000	ng/ul	-0.04
64) Phenanthrene-d10	17.098	188		751688	20.000	ng/ul	-0.02
79) Chrysene-d12	21.545	240		859798	20.000	ng/ul	# 0.00
88) Perylene-d12	24.845	264		1003259	20.000	ng/ul	0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.187	96		17779	6.158	ng/uL	-0.04
4) Pyridine-d5	3.593	84		246685	29.889	ng/ul	-0.04
7) Phenol-d5	6.822	99		280002	30.266	ng/ul	-0.07
9) Bis-(2-Chloroethyl)eth...	6.975	67		165408	29.036	ng/ul	-0.07
11) 2-Chlorophenol-d4	7.181	132		213580	31.909	ng/ul	-0.06
15) 4-Methylphenol-d8	8.334	113		220786	30.118	ng/ul	-0.07
21) Nitrobenzene-d5	8.781	128		106586	33.948	ng/ul	-0.06
24) 2-Nitrophenol-d4	9.498	143		125099	35.141	ng/ul	-0.06
28) 2,4-Dichlorophenol-d3	10.028	165		238631	30.867	ng/ul	-0.06
31) 4-Chloroaniline-d4	10.545	131		286600	31.070	ng/ul	-0.06
46) Dimethylphthalate-d6	13.698	166		837454	34.250	ng/ul	-0.02
49) Acenaphthylene-d8	13.975	160		877696	33.904	ng/ul	-0.03
54) 4-Nitrophenol-d4	14.504	143		117382	28.066	ng/ul	-0.02
60) Fluorene-d10	15.304	176		715476	33.080	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.439	200		135611	29.059	ng/ul	0.00
73) Anthracene-d10	17.204	188		1204162	34.315	ng/ul	-0.01
81) Pyrene-d10	19.586	212		1547514	34.741	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.615	264		1783886	33.806	ng/ul	0.00

Target Compounds Qvalue

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

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