

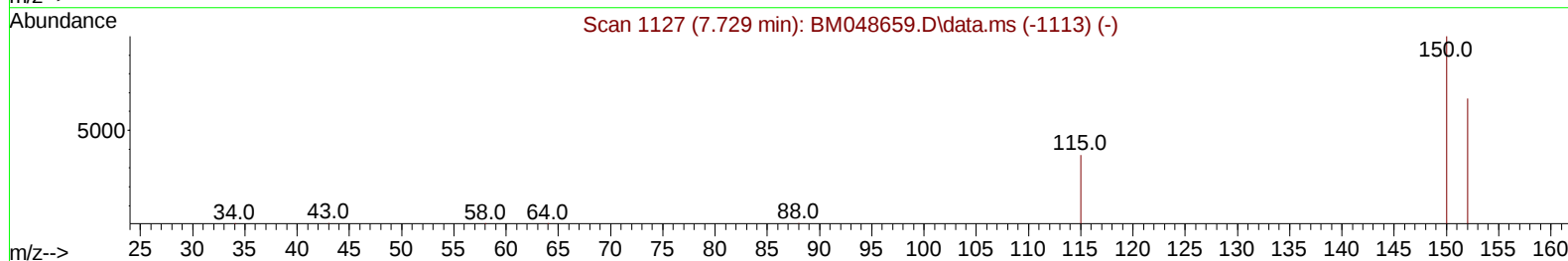
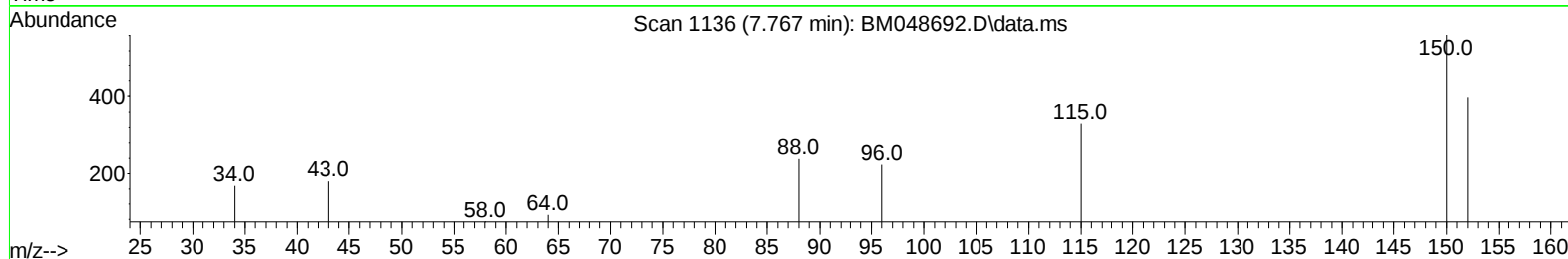
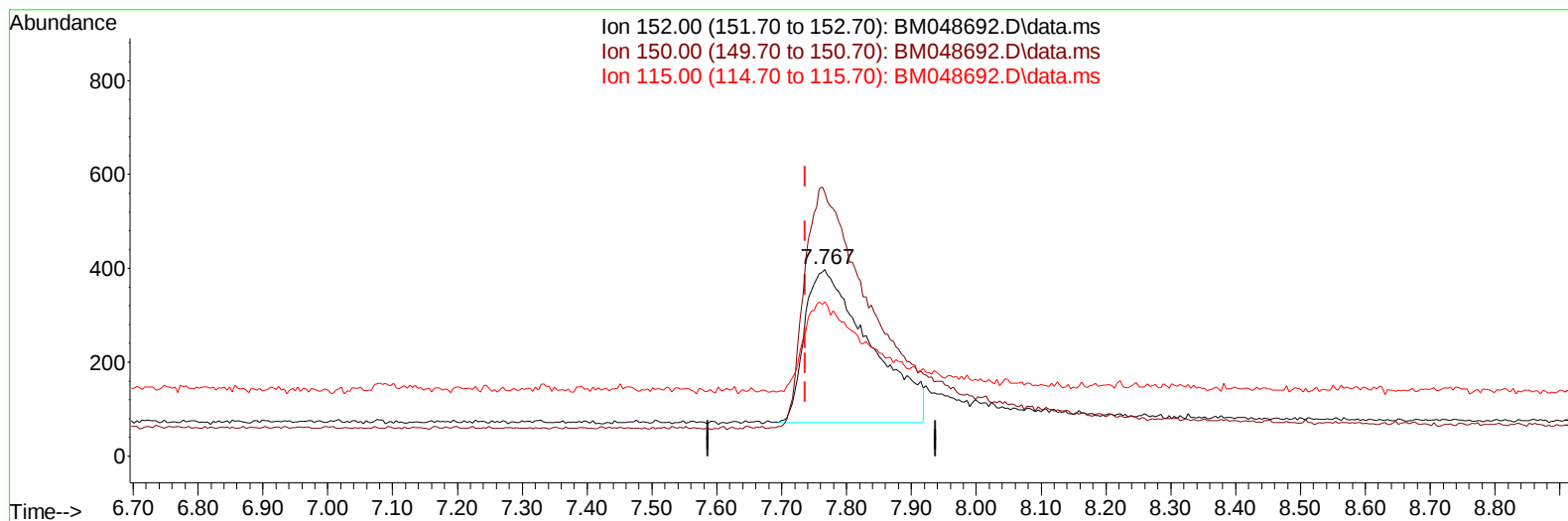
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111324\
Data File : BM048692.D
Acq On : 15 Nov 2024 02:58
Operator : RC/JU
Sample : PB164925BS
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SLCS925

Manual IntegrationsAPPROVED

Quant Time: Nov 15 03:31:26 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 13 11:38:55 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/15/2024



TIC: BM048692.D\data.ms

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

7.767min (+ 0.030) 0.40 ng/u1

response 2272

Ion	Exp%	Act%
152.00	100.00	100.00
150.00	151.00	140.95
115.00	65.50	82.66#
0.00	0.00	0.00

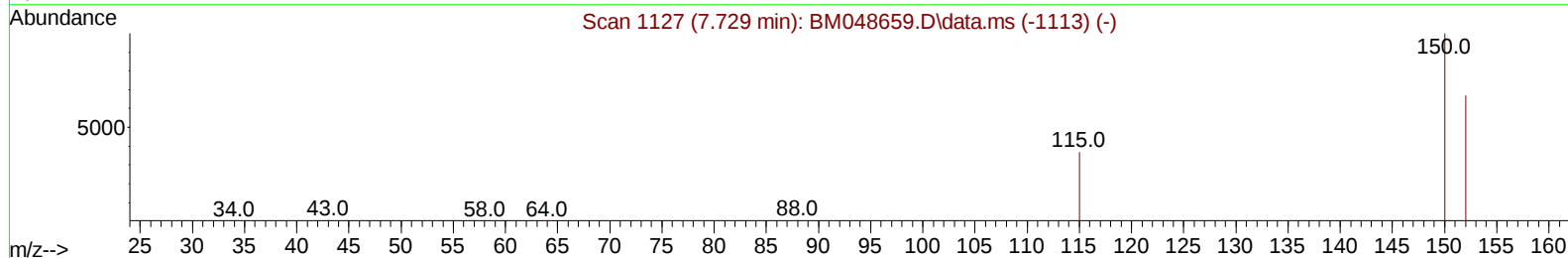
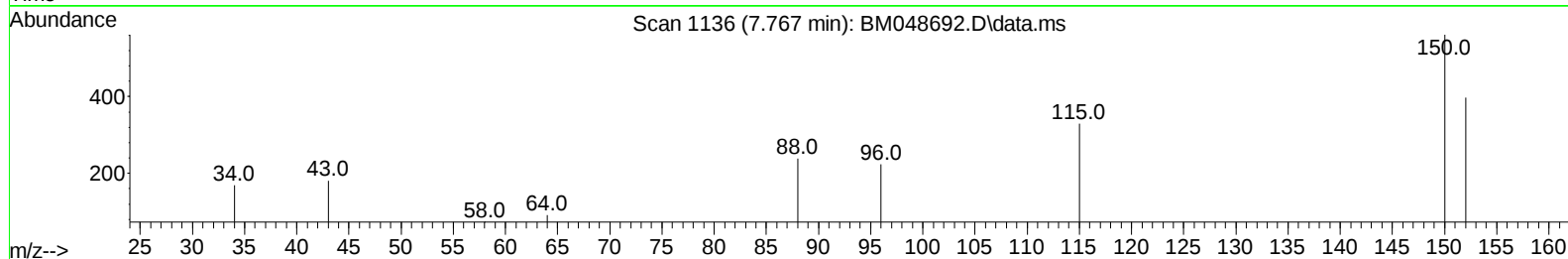
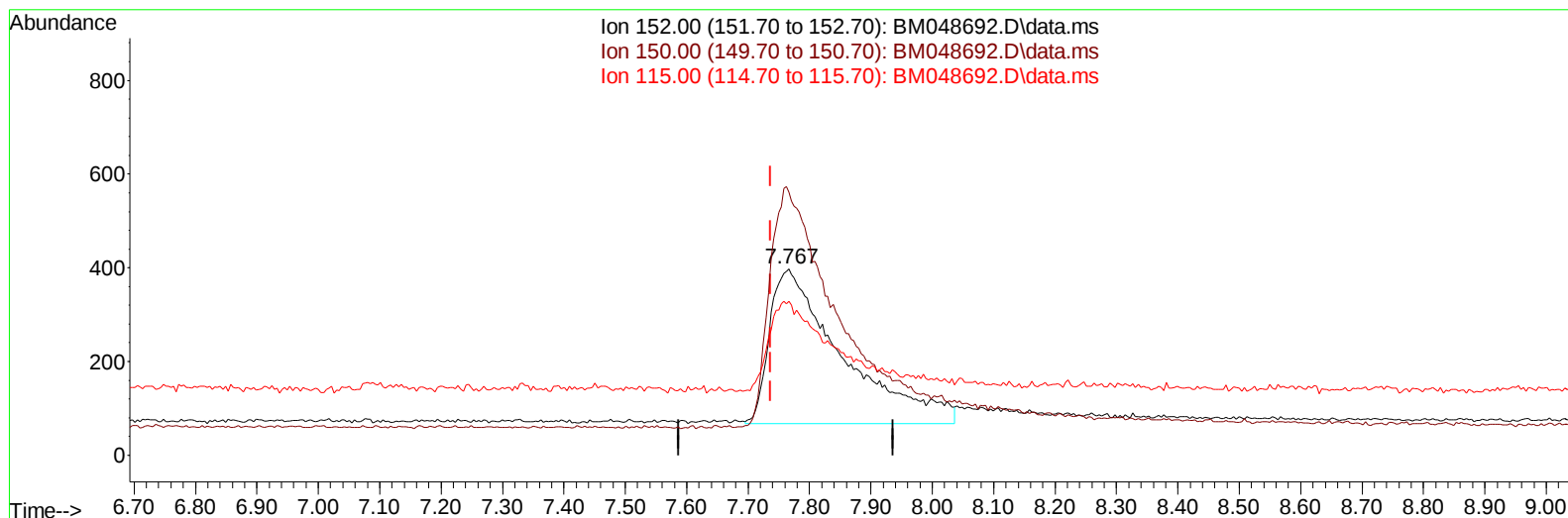
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(1) 1,4-Dichlorobenzene-d4 (I)

7.767min (+ 0.030) 0.40 ng/ul m

response 2698

Ion	Exp%	Act%
152.00	100.00	100.00
150.00	151.00	140.95
115.00	65.50	82.66#
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.767	152		2698m	0.400	ng/ul	0.03
4) Naphthalene-d8	10.506	136		9066	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.325	164		5495	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.089	188		11135	0.400	ng/ul	0.00
17) Chrysene-d12	21.266	240		8990	0.400	ng/ul	0.00
23) Perylene-d12	23.484	264		9140	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.122	96		3000	0.922	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.112	152		4975	0.424	ng/ul	0.00
18) Fluoranthene-d10	19.112	212		11638	0.461	ng/ul	0.00
Target Compounds							
						Qvalue	
2) 1,4-Dioxane	3.155	88		8351	2.232	ng/ul#	63
5) Naphthalene	10.556	128		9743	0.429	ng/ul	96
7) 2-Methylnaphthalene	12.183	142		5590	0.391	ng/ul	94
8) 1-Methylnaphthalene	12.381	142		6689	0.457	ng/ul	99
10) Acenaphthylene	14.052	152		9098	0.436	ng/ul	98
11) Acenaphthene	14.385	153		7268	0.435	ng/ul	99
12) Fluorene	15.403	166		7712	0.425	ng/ul	100
14) Pentachlorophenol	16.751	266		2428	0.594	ng/ul	99
15) Phenanthrene	17.131	178		11320	0.387	ng/ul	98
16) Anthracene	17.245	178		10110	0.371	ng/ul	97
19) Fluoranthene	19.140	202		15170	0.456	ng/ul	99
20) Pyrene	19.498	202		16354	0.453	ng/ul	99
21) Benzo(a)anthracene	21.254	228		10642	0.362	ng/ul	99
22) Chrysene	21.298	228		18068	0.512	ng/ul	100
24) Benzo(b)fluoranthene	22.815	252		13571	0.476	ng/ul	96
25) Benzo(k)fluoranthene	22.861	252		17638	0.528	ng/ul	96
26) Benzo(a)pyrene	23.396	252		13310	0.479	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.749	276		16185	0.417	ng/ul#	50
28) Dibenzo(a,h)anthracene	25.769	278		11812	0.391	ng/ul	96
29) Benzo(g,h,i)perylene	26.420	276		14738	0.452	ng/ul	99

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

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