

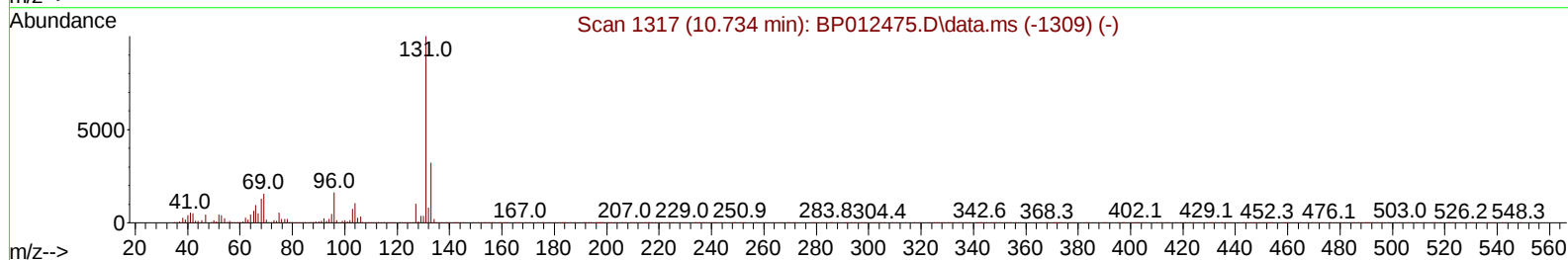
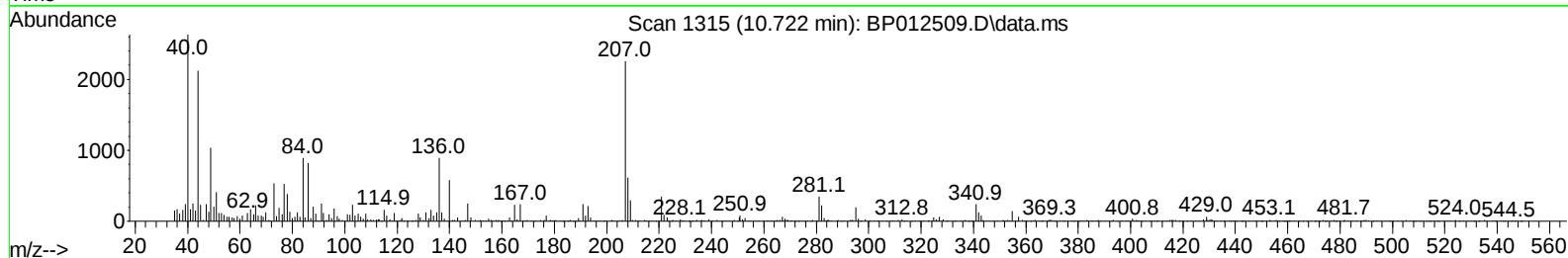
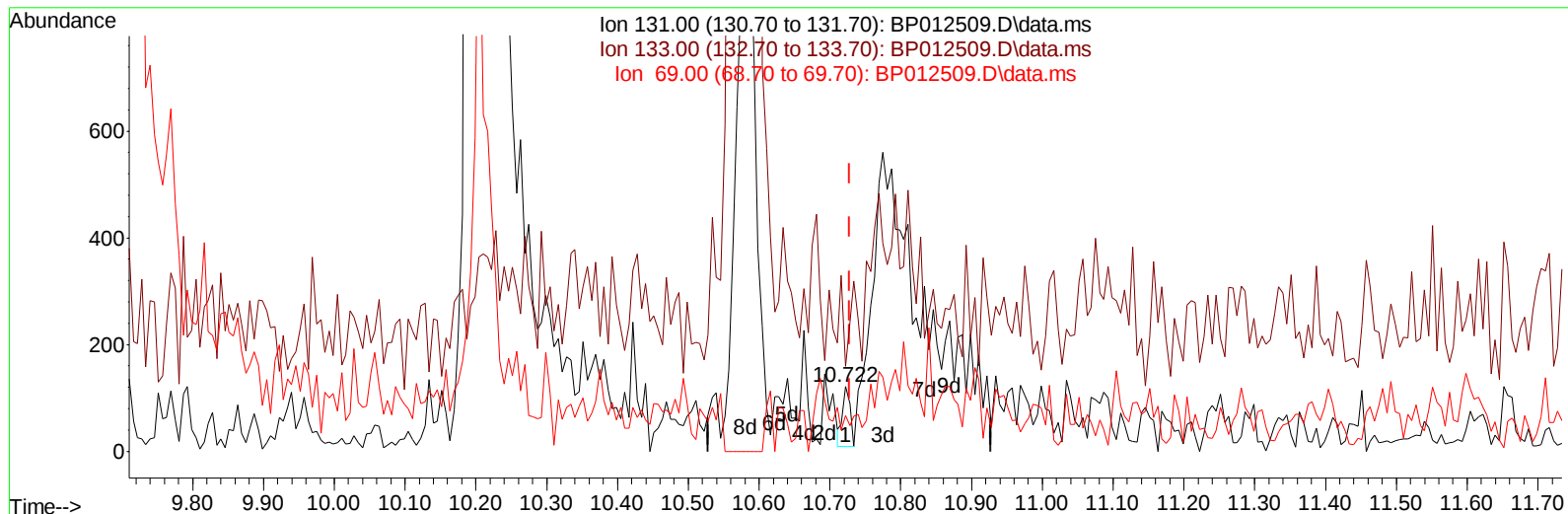
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012509.D
 Acq On : 04 Nov 2022 03:01
 Operator : CG/JU
 Sample : PB148748BL
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK748

Manual IntegrationsAPPROVED

Quant Time: Nov 04 03:53:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/05/2022
 Supervised By :mohammad ahmed 11/06/2022



TIC: BP012509.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.722min (-0.006) 0.00 ng/u1

response 83

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	131.40#
69.00	16.50	55.37#
0.00	0.00	0.00

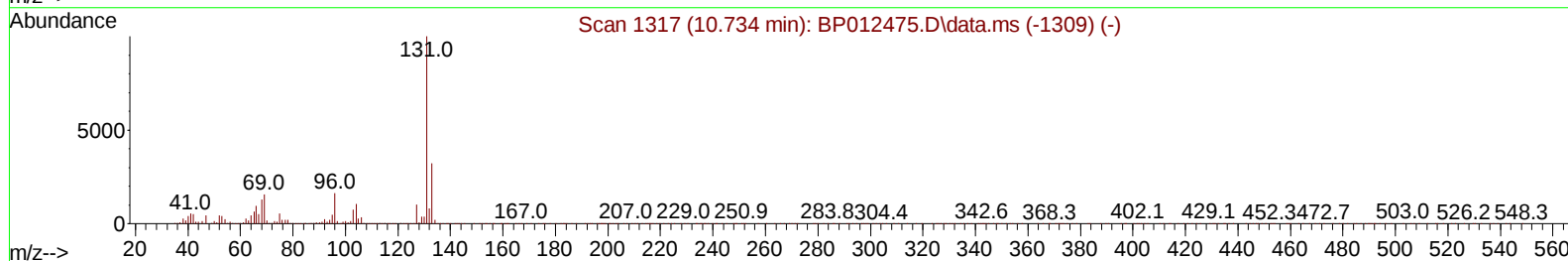
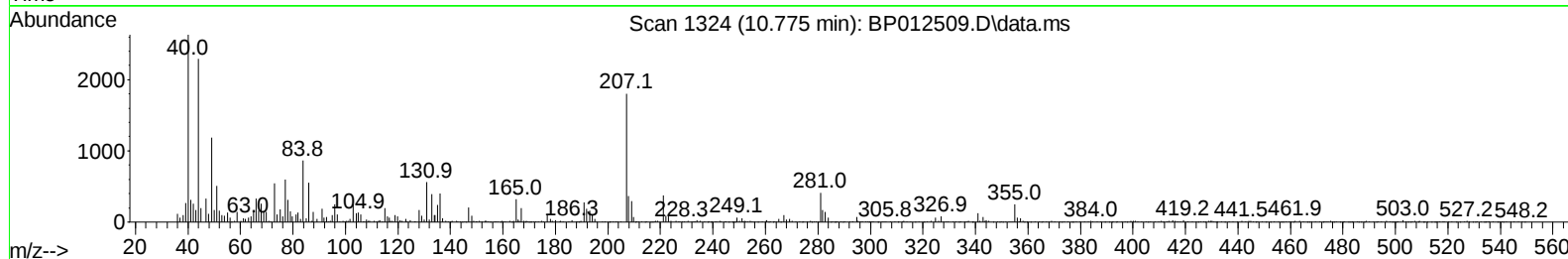
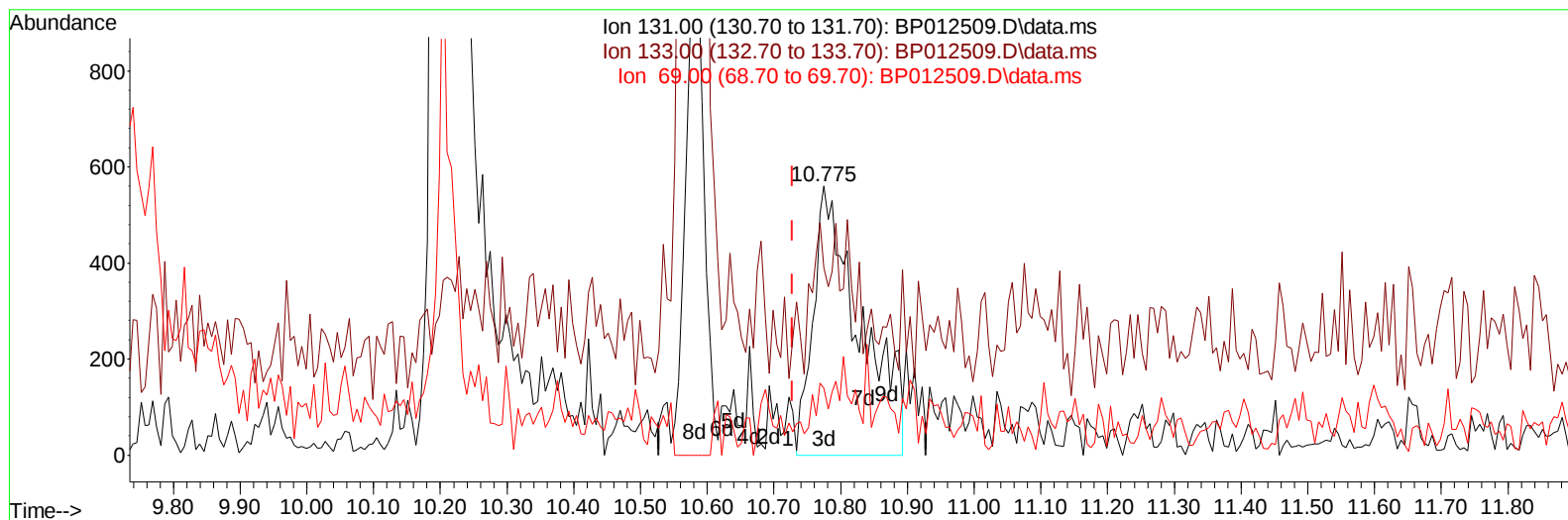
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TIC: BP012509.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.775min (+ 0.047) 0.10 ng/ul m

response 2724

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	69.64#
69.00	16.50	24.82#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	259643	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1236062	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	873890	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	1959754	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1678882	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1490495	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	38170	5.919	ng/uL	0.00
4) Pyridine-d5	3.669	84	1020	0.055	ng/ul	0.02
7) Phenol-d5	6.952	99	746198	32.042	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	431356	31.028	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	570945	33.572	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	609862	33.111	ng/ul	0.00
21) Nitrobenzene-d5	8.946	128	297200	37.078	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	338647	40.704	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	589000	32.667	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	2724m	0.101	ng/ul	0.05
46) Dimethylphthalate-d6	13.845	166	2034167	34.003	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	2186936	31.990	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	286318	25.288	ng/ul	0.00
60) Fluorene-d10	15.428	176	1741336	33.998	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	205226	20.114	ng/ul	0.00
73) Anthracene-d10	17.292	188	2844462	33.686	ng/ul	0.00
81) Pyrene-d10	19.533	212	3164340	37.541	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	2482306	34.085	ng/ul	0.00

Target Compounds Qvalue

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

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