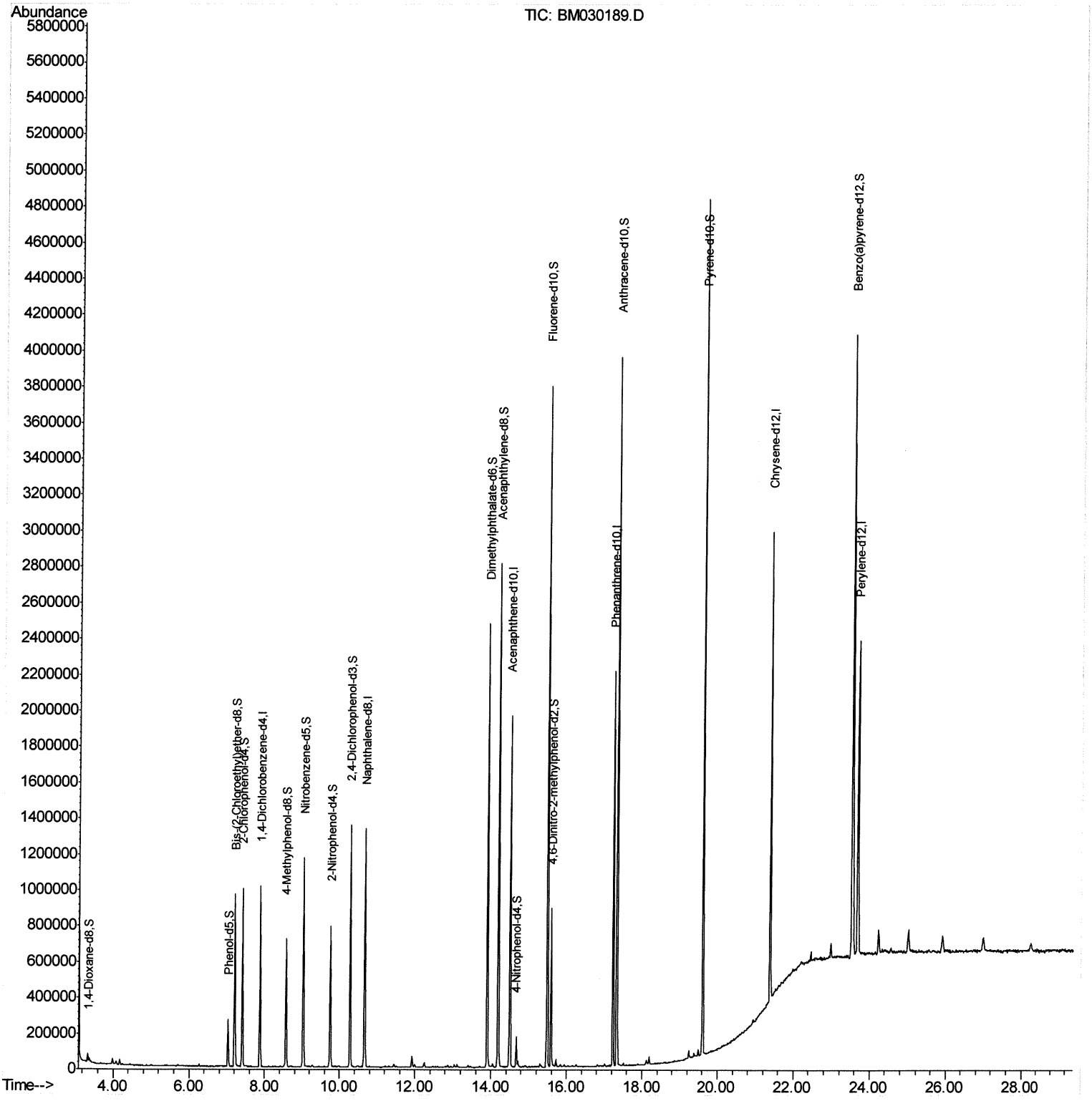
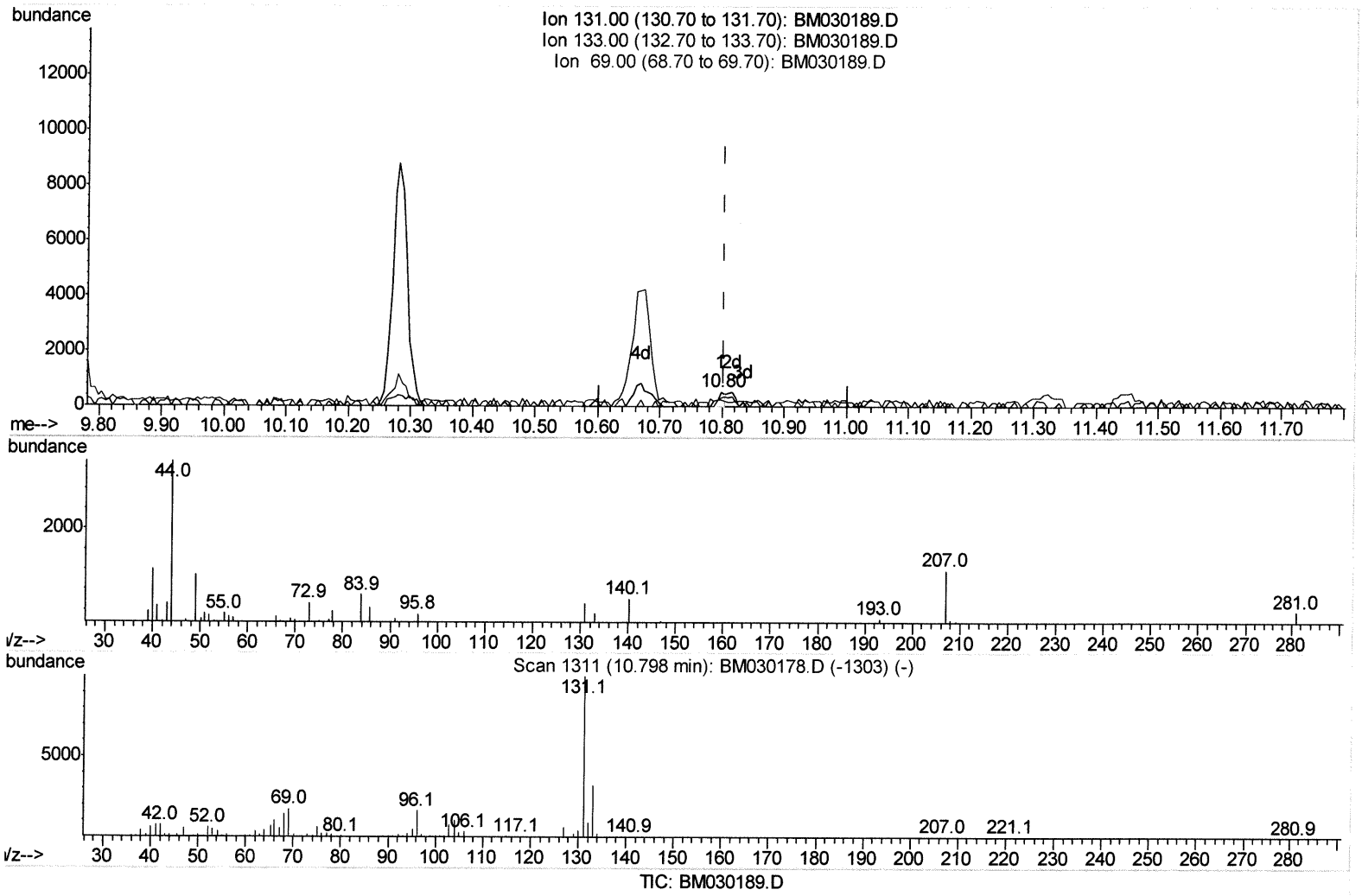
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052621\
Data File : BM030189.D
Acq On : 27 May 2021 03:58
Operator : CG/JU
Sample : M2495-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: May 27 04:40:50 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM052621.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 26 15:30:24 2021
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052621\
Data File : BM030189.D
Acq On : 27 May 2021 03:58
Operator : CG/JU
Sample : M2495-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: May 27 04:39:16 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM052621.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 26 15:30:24 2021
Response via : Initial Calibration



TIC: BM030189.D

(31) 4-Chloroaniline-d4 (S)

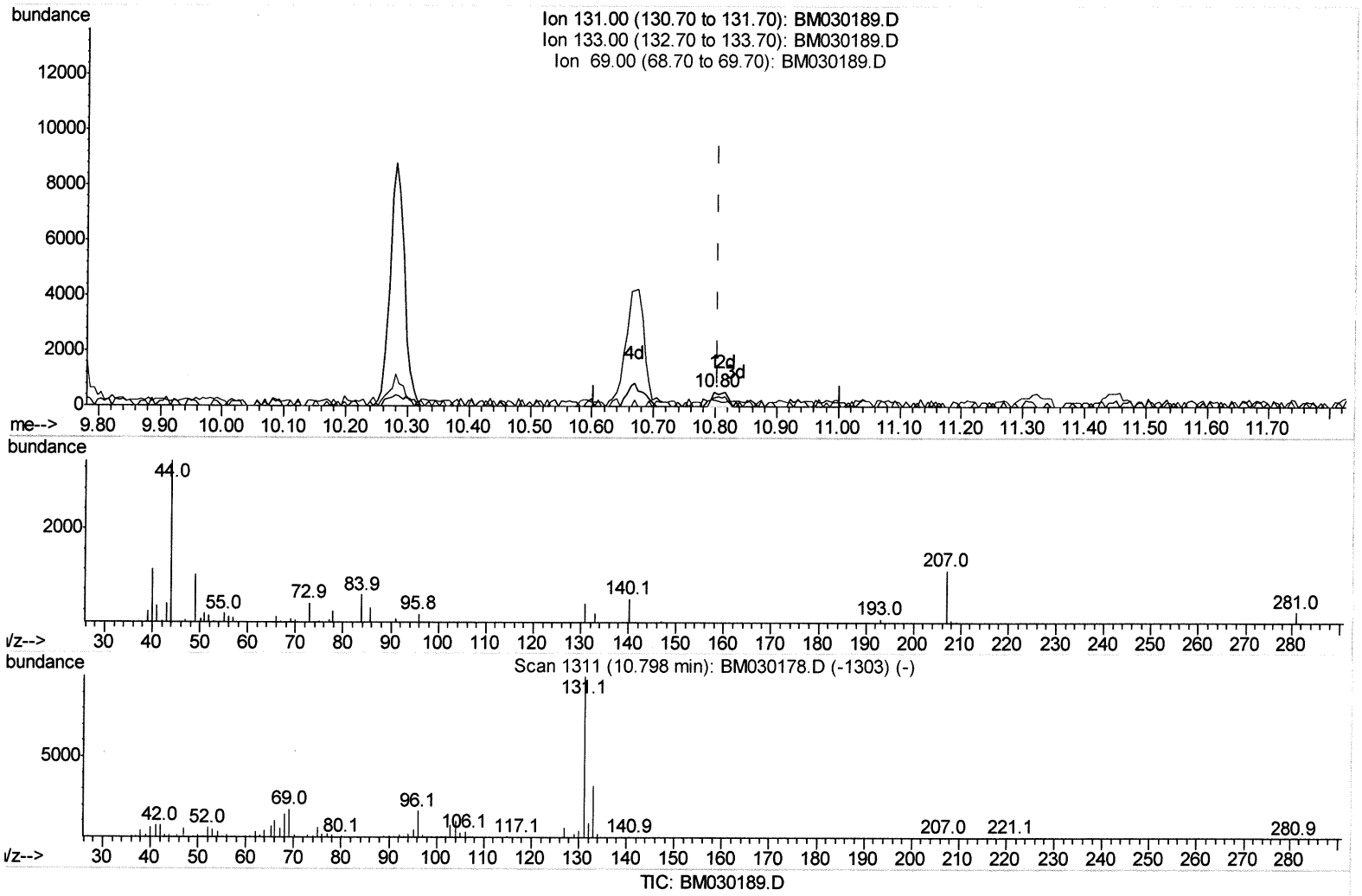
10.798min (-0.005) 0.02ng/ul

response 498

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	63.03#
69.00	17.50	42.91#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052621\
Data File : BM030189.D
Acq On : 27 May 2021 03:58
Operator : CG/JU
Sample : M2495-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: May 27 04:39:16 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM052621.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 26 15:30:24 2021
Response via : Initial Calibration



TIC: BM030189.D

(31) 4-Chloroaniline-d4 (S)

10.798min (-0.005) 0.04ng/ul m JU 5/28/21

response 956

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	63.03#
69.00	17.50	42.91#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052621\
Data File : BM030189.D
Acq On : 27 May 2021 03:58
Operator : CG/JU
Sample : M2495-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: May 27 04:40:50 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM052621.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 26 15:30:24 2021
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	269216	20.00	ng/ul	0.00
20) Naphthalene-d8	10.67	136	1087343	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.50	164	671829	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.23	188	1292971	20.00	ng/ul	0.00
79) Chrysene-d12	21.39	240	1189474	20.00	ng/ul	0.00
88) Perylene-d12	23.68	264	1212688	20.00	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.34	96	22871	3.48	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	7.03	99	134280	5.67	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.20	67	434516	29.08	ng/ul	0.00
11) 2-Chlorophenol-d4	7.40	132	430850	24.30	ng/ul	0.00
15) 4-Methylphenol-d8	8.57	113	253575	13.49	ng/ul	0.00
21) Nitrobenzene-d5	9.03	128	250661	35.32	ng/ul	0.00
24) 2-Nitrophenol-d4	9.75	143	241094	35.67	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.28	165	482548	27.67	ng/ul	0.00
31) 4-Chloroaniline-d4	10.80	131	956m	0.04	ng/ul	0.00
46) Dimethylphthalate-d6	13.90	166	1574031	30.65	ng/ul	0.00
49) Acenaphthylene-d8	14.19	160	1947860	29.42	ng/ul	0.00
54) 4-Nitrophenol-d4	14.67	143	35851	4.32	ng/ul	-0.01
60) Fluorene-d10	15.49	176	1401956	31.27	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.59	200	179957	28.71	ng/ul	-0.01
73) Anthracene-d10	17.33	188	2144178	33.51	ng/ul	0.00
81) Pyrene-d10	19.61	212	2293435	35.97	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.53	264	2284832	33.24	ng/ul	-0.01

Target Compounds	Qvalue

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

JH 5/28/21