

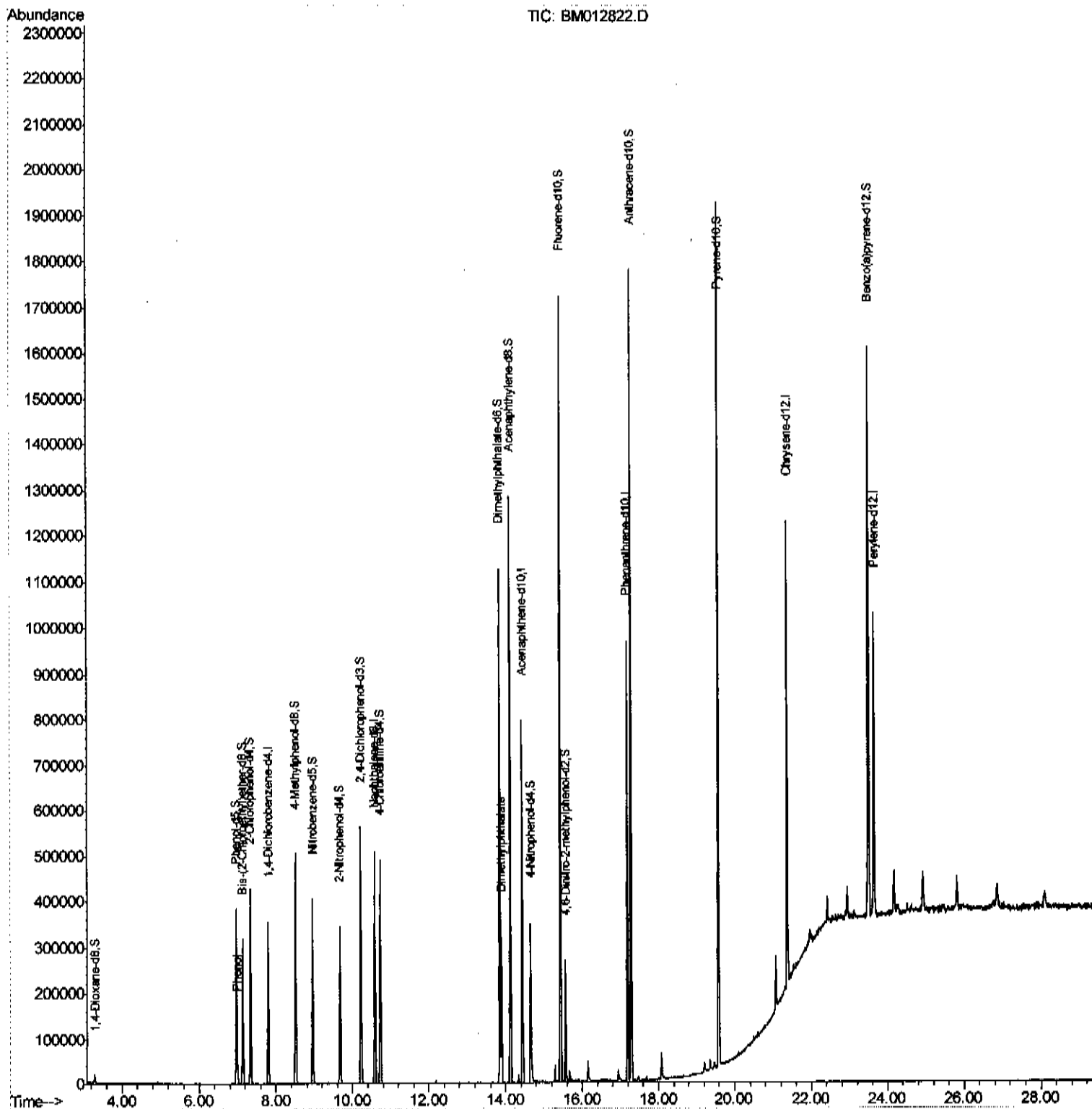
Data Path : Z:\HPCHEM1\BNA M\DATA\BM110817\
Data File : BM012822.D
Acq On : 08 Nov 2017 15:45
Operator : SJ/JU
Sample : I5955-11
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-14(7-7.5)-101717

Manual Integrations
APPROVED

Sohil
11/9/2017 3:48:28 PM

Quant Time: Nov 09 03:27:14 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 09 01:31:40 2017
Response via : Initial Calibration



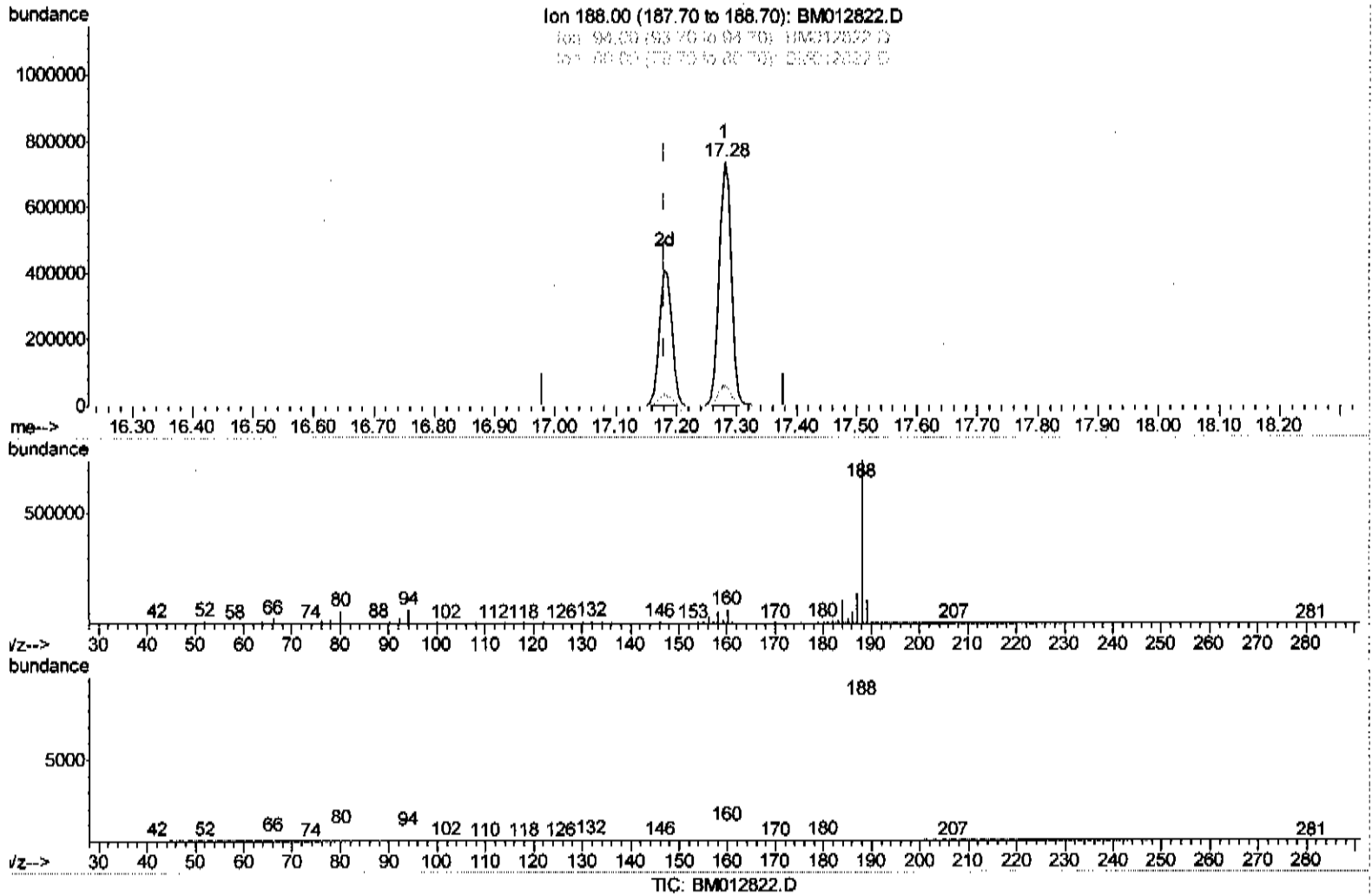
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(61) Phenanthrene-d10 (I)

17.280min (+0.100) 20.00ng/ul

response 1021942

Ion	Exp%	Act%
188.00	100	100
94.00	9.00	8.44
80.00	9.70	7.78
0.00	0.00	0.00

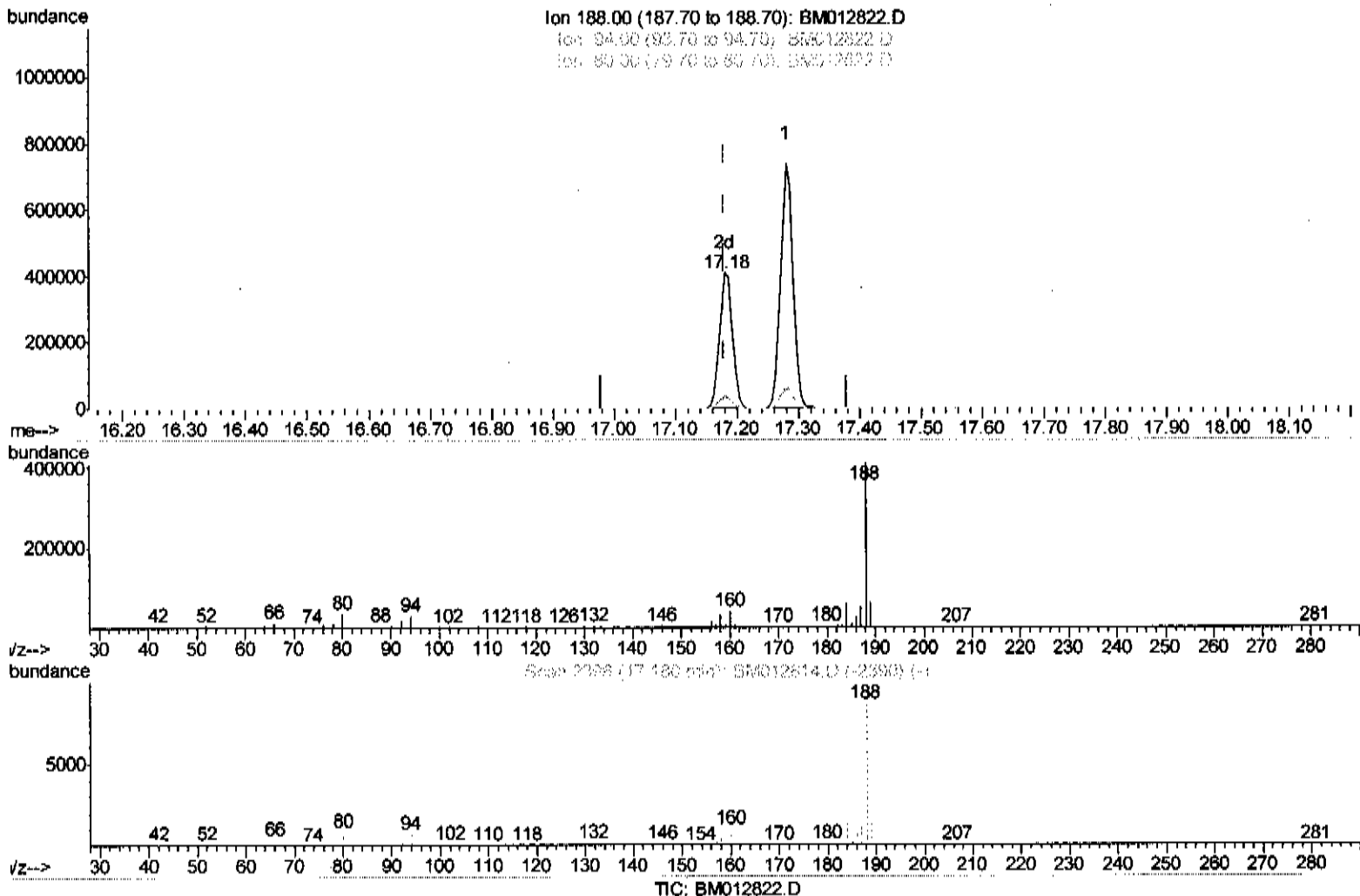
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(61) Phenanthrene-d10 (I)

17.180min (+0.000) 20.00ng/ul m

Ju 11/13/17

response 591342

Ion	Exp%	Act%
188.00	100	100
94.00	9.00	7.15#
80.00	9.70	8.71
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	101223	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	426318	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	269375	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.18	188	591342m	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	485552	20.00	ng/ul	0.00
85) Perylene-d12	23.64	264	508657	20.00	ng/ul	0.00

10/11/13/17

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	10036	5.36	ng/uL	0.00
5) Phenol-d5	6.97	99	231493	30.74	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	130544	31.98	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	197940	31.55	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	203192	31.25	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	102363	32.50	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	119514	33.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	224427	30.37	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	274251	38.46	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	780883	35.03	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	929797	33.72	ng/ul	0.00
51) 4-Nitrophenol-d4	14.65	143	104697	29.31	ng/ul	0.00
57) Fluorene-d10	15.43	176	659064	34.52	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.57	200	72045	18.45	ng/ul	0.00
70) Anthracene-d10	17.28	188	1021942	35.77	ng/ul	0.00
78) Pyrene-d10	19.57	212	976560	43.34	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.50	264	893099	37.13	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.00	94	19793	2.56	ng/ul	92
44) Dimethylphthalate	13.89	163	208383	9.42	ng/ul	98

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