

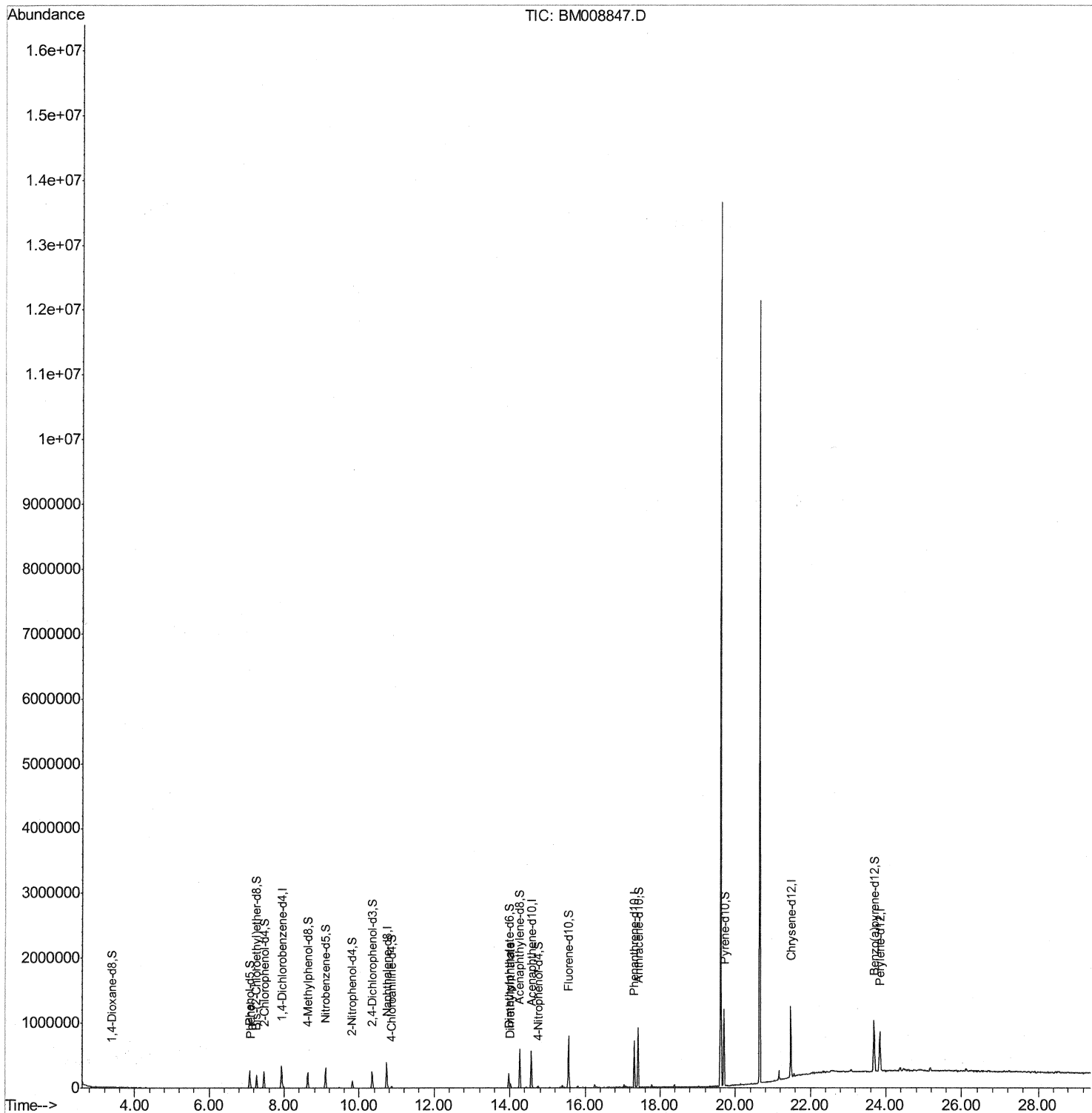
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012417\
 Data File : BM008847.D
 Acq On : 24 Jan 2017 20:27
 Operator : UM/SJ
 Sample : I1315-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDN28

Manual Integrations
 APPROVED

mohammad
 1/25/2017 3:46:11 PM

Quant Time: Jan 25 03:51:02 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 25 03:34:06 2017
 Response via : Initial Calibration



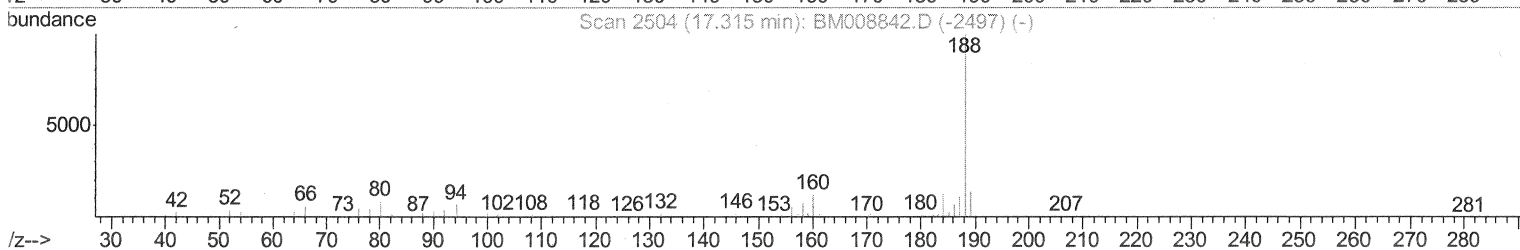
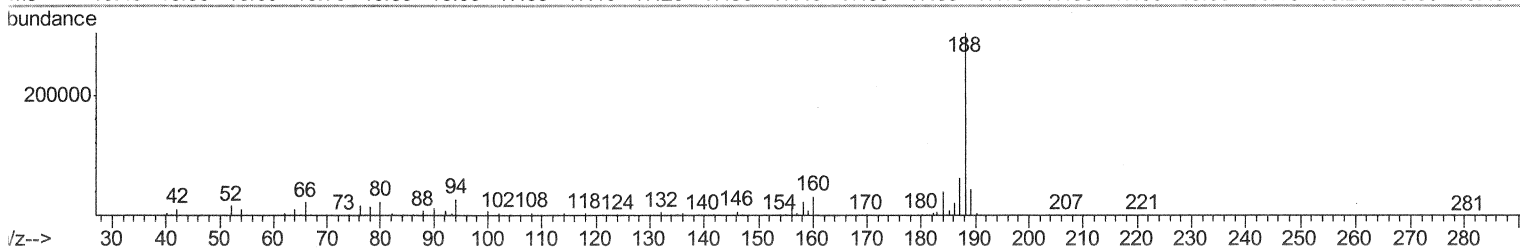
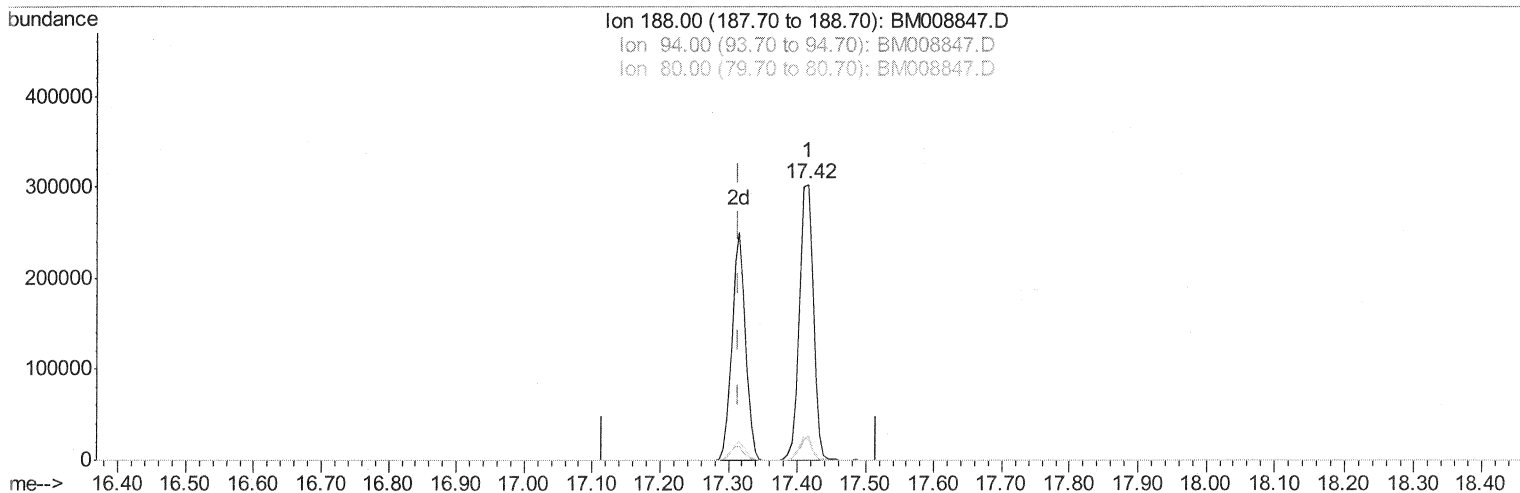
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TIC: BM008847.D

(61) Phenanthrene-d10 (I)

17.415min (+0.100) 20.00ng/ul

response 428843

Ion	Exp%	Act%
188.00	100	100
94.00	8.00	9.08
80.00	9.50	7.76
0.00	0.00	0.00

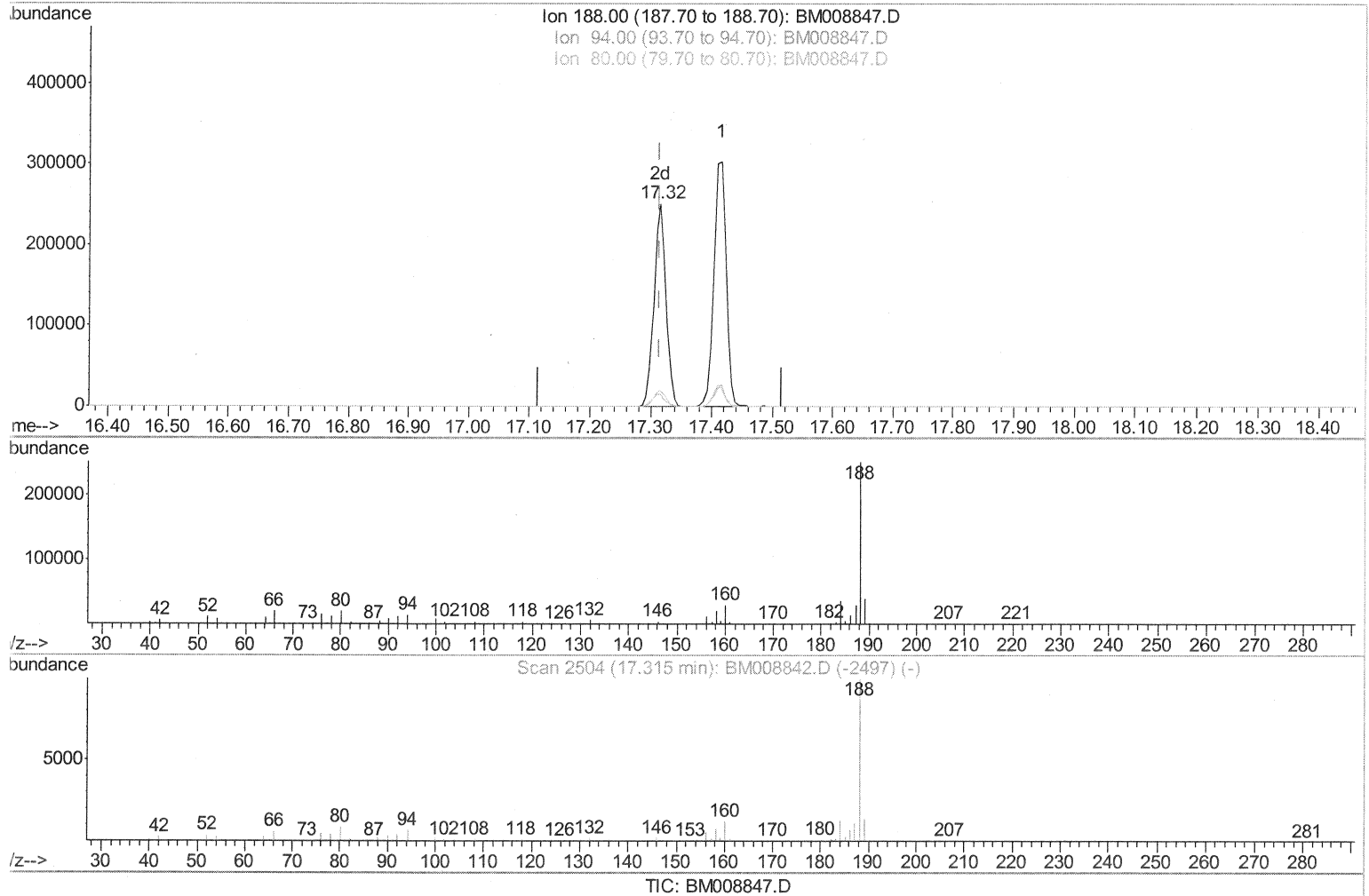
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(61) Phenanthrene-d10 (I)

17.315min (-0.000) 20.00ng/ul m SJ 1/12/17

response 348756

Ion	Exp%	Act%
188.00	100	100
94.00	8.00	5.86#
80.00	9.50	8.20
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	52916	20.00	ng/ul	0.00
18) Naphthalene-d8	10.74	136	208508	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.57	164	157482	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.32	188	348756m>	20.00	ng/ul	0.00
75) Chrysene-d12	21.48	240	483079	20.00	ng/ul	0.00
83) Perylene-d12	23.84	264	453437	20.00	ng/ul	0.00

SJ 1/27/17

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	1234	1.15	ng/uL	0.01
5) Phenol-d5	7.08	99	81064	18.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.26	67	89528	25.13	ng/ul	0.00
9) 2-Chlorophenol-d4	7.46	132	52638	18.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.63	113	54925	15.65	ng/ul	0.00
19) Nitrobenzene-d5	9.10	128	31213	23.54	ng/ul	0.00
22) 2-Nitrophenol-d4	9.82	143	16964	10.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.35	165	55469	16.43	ng/ul	0.00
29) 4-Chloroaniline-d4	10.88	131	6452	1.76	ng/ul	0.00
43) Dimethylphthalate-d6	13.97	166	96378	9.34	ng/ul	0.00
46) Acenaphthylene-d8	14.26	160	303372	24.85	ng/ul	0.00
51) 4-Nitrophenol-d4	14.74	143	3741	2.17	ng/ul	-0.01
57) Fluorene-d10	15.56	176	246685	25.17	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.67	200	606	0.29	ng/ul	0.00
70) Anthracene-d10	17.42	188	428843	28.12	ng/ul	0.00
76) Pyrene-d10	19.70	212	544374	25.69	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.70	264	502231	26.15	ng/ul	0.00

Target Compounds

						Qvalue
6) Phenol	7.11	94	10770	2.20	ng/ul#	48
44) Dimethylphthalate	14.02	163	24978	2.42	ng/ul#	95

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