

Data Path : Z:\HPCHEM1\BNA_M\Data\BM031117\
 Data File : BM009211.D
 Acq On : 11 Mar 2017 17:06
 Operator : SJ/MA
 Sample : I2107-17DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AL3DL

Quant Time: Mar 13 12:04:27 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 11:21:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	228	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	702	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	392	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	906	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	920	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	916	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	39	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	104	0.04	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.73	128	41	0.02	ng/ul#	1
7) Acenaphthylene	14.23	152	49	0.02	ng/ul#	1
8) Acenaphthene	14.57	153	32	0.02	ng/ul#	77
9) Fluorene	15.56	166	42	0.02	ng/ul#	62
12) Phenanthrene	17.30	178	686	0.24	ng/ul#	85
13) Anthracene	17.40	178	162	0.06	ng/ul#	33
15) Fluoranthene	19.31	202	1255	0.34	ng/ul	95
17) Pyrene	19.68	202	1057	0.33	ng/ul#	89
18) Benzo(a)anthracene	21.42	228	520	0.17	ng/ul#	83
19) Chrysene	21.47	228	504	0.17	ng/ul#	79
21) Benzo(b)fluoranthene	23.07	252	916	0.25	ng/ul	69
22) Benzo(k)fluoranthene	23.07	252	916	0.29	ng/ul#	67
23) Benzo(a)pyrene	23.67	252	490	0.15	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	26.15	276	338	0.08	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.15	278	91	0.03	ng/ul#	1
26) Benzo(g,h,i)perylene	26.88	276	313	0.09	ng/ul#	57

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

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