

Data Path : Z:\HPCHEM1\BNA_M\Data\BM031117\
 Data File : BM009212.D
 Acq On : 11 Mar 2017 17:42
 Operator : SJ/MA
 Sample : I2072-06DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK1DL

Quant Time: Mar 13 12:07:33 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	203	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	655	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	373	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	889	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1014	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1089	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.27	152	94	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	247	0.09	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.73	128	168	0.09	ng/ul#	38
5) 2-Methylnaphthalene	12.34	142	108	0.08	ng/ul	95
7) Acenaphthylene	14.23	152	39	0.02	ng/ul#	1
12) Phenanthrene	17.30	178	1327	0.48	ng/ul	94
13) Anthracene	17.30	178	1315	0.48	ng/ul	95
15) Fluoranthene	19.31	202	1834	0.50	ng/ul	98
17) Pyrene	19.69	202	2390	0.69	ng/ul	96
18) Benzo(a)anthracene	21.42	228	953	0.28	ng/ul#	90
19) Chrysene	21.42	228	953	0.30	ng/ul	92
21) Benzo(b)fluoranthene	23.07	252	1866	0.44	ng/ul	83
22) Benzo(k)fluoranthene	23.07	252	1851	0.49	ng/ul#	81
23) Benzo(a)pyrene	23.67	252	717	0.19	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.15	276	646	0.13	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.15	278	176	0.04	ng/ul#	6
26) Benzo(g,h,i)perylene	26.88	276	563	0.13	ng/ul#	76

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

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