

Data Path : Z:\HPCHEM1\BNA_M\Data\BM031117\
 Data File : BM009207.D
 Acq On : 11 Mar 2017 14:43
 Operator : SJ/MA
 Sample : I2107-04DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ3DL

Quant Time: Mar 13 11:47:23 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	243	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	792	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	454	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1054	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1214	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	1218	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	27	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	81	0.02	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.73	128	392	0.17	ng/ul#	69
5) 2-Methylnaphthalene	12.34	142	159	0.10	ng/ul	99
7) Acenaphthylene	14.23	152	117	0.05	ng/ul#	15
8) Acenaphthene	14.57	153	58	0.03	ng/ul#	82
9) Fluorene	15.56	166	78	0.04	ng/ul#	62
12) Phenanthrene	17.30	178	1446	0.44	ng/ul	94
13) Anthracene	17.30	178	1438	0.44	ng/ul	95
15) Fluoranthene	19.31	202	3129	0.72	ng/ul	97
17) Pyrene	19.68	202	2662	0.64	ng/ul	96
18) Benzo(a)anthracene	21.42	228	1732	0.43	ng/ul	92
19) Chrysene	21.47	228	1814	0.47	ng/ul	91
21) Benzo(b)fluoranthene	23.07	252	3955	0.83	ng/ul	97
22) Benzo(k)fluoranthene	23.07	252	3955	0.93	ng/ul	96
23) Benzo(a)pyrene	23.57	252	1430	0.33	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.15	276	1505	0.27	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.16	278	377	0.08	ng/ul#	43
26) Benzo(g,h,i)perylene	26.88	276	1405	0.30	ng/ul	94

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

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