

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061815\
 Data File : BM001734.D
 Acq On : 18 Jun 2015 11:06
 Operator : TP/IZ
 Sample : G2610-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1W49

Quant Time: Jun 18 17:15:13 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 18 04:33:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	918	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	3275	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1550	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3246	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	2642	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	2326	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1364	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	2470	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.50	128	4013	0.46	ng/ul	100
5) 2-Methylnaphthalene	12.12	142	5127	0.93	ng/ul	100
7) Acenaphthylene	14.03	152	4214	0.56	ng/ul	98
8) Acenaphthene	14.37	153	7739	1.36	ng/ul	98
11) Pentachlorophenol	16.70	266	437	0.63	ng/ul	100
12) Phenanthrene	17.09	178	6109	0.63	ng/ul#	90
15) Fluoranthene	19.12	202	10600	1.02	ng/ul	86
17) Pyrene	19.48	202	8559	0.81	ng/ul	92
18) Benzo(a)anthracene	21.23	228	4431	0.52	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	7019	0.78	ng/ul	96
22) Benzo(k)fluoranthene	22.85	252	8566	0.99	ng/ul	96
23) Benzo(a)pyrene	23.37	252	4124	0.51	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.70	276	7398	0.79	ng/ul	97
25) Dibenzo(a,h)anthracene	25.73	278	7379	0.97	ng/ul	95
26) Benzo(g,h,i)perylene	26.39	276	8314	1.02	ng/ul	97

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

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