

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072816\
 Data File : BM006726.D
 Acq On : 28 Jul 2016 17:51
 Operator : UM/SJ
 Sample : H4068-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2548

Quant Time: Jul 28 18:35:44 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM072816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 15:44:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	2176	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7747	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4168	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	10270	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	10293	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	10042	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	3496	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	9842	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	12160	0.62	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	5481	0.41	ng/ul	100
9) Fluorene	15.32	166	9962	0.59	ng/ul	95
11) Pentachlorophenol	16.68	266	4998	3.87	ng/ul	96
13) Anthracene	17.14	178	21971	0.81	ng/ul	98
15) Fluoranthene	19.08	202	24733	0.62	ng/ul	94
17) Pyrene	19.45	202	26978	0.87	ng/ul	96
19) Chrysene	21.25	228	18783	0.66	ng/ul	93
21) Benzo(b)fluoranthene	22.76	252	17198	0.50	ng/ul	94
23) Benzo(a)pyrene	23.33	252	21957	0.69	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.63	278	44662	1.49	ng/ul#	95

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

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