

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072816\
 Data File : BM006727.D
 Acq On : 28 Jul 2016 18:28
 Operator : UM/SJ
 Sample : H4068-03DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2548DL

Quant Time: Jul 28 19:01:42 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	2111	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7334	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3960	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	9751	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	9998	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	9912	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1676	0.18	ng/ul	0.01
14) Fluoranthene-d10	19.05	212	5179	0.20	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	5957	0.32	ng/ul	95
5) 2-Methylnaphthalene	12.07	142	2627	0.21	ng/ul	99
9) Fluorene	15.32	166	4854	0.30	ng/ul	90
11) Pentachlorophenol	16.70	266	2057	1.68	ng/ul	95
13) Anthracene	17.14	178	10564	0.41	ng/ul	96
15) Fluoranthene	19.08	202	12602	0.33	ng/ul	95
17) Pyrene	19.45	202	13344	0.44	ng/ul	94
19) Chrysene	21.26	228	9761	0.36	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	8938	0.26	ng/ul	96
23) Benzo(a)pyrene	23.32	252	11083	0.35	ng/ul	98
25) Dibenzo(a,h)anthracene	25.62	278	22807	0.77	ng/ul	96

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

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