

Data Path : Z:\HPCHEM1\BNA_M\Data\BM072916\
 Data File : BM006733.D
 Acq On : 29 Jul 2016 10:56
 Operator : UM/SJ
 Sample : H4068-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2553

Quant Time: Jul 29 12:43:06 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	1731	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6115	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3295	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	8014	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8907	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	8465	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	2447	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	7891	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	8832	0.57	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	5553	0.53	ng/ul	99
9) Fluorene	15.32	166	8679	0.65	ng/ul	99
11) Pentachlorophenol	16.68	266	1833	1.82	ng/ul	96
12) Phenanthrene	17.06	178	15877	0.71	ng/ul	98
15) Fluoranthene	19.08	202	17943	0.58	ng/ul	97
18) Benzo(a)anthracene	21.20	228	20206	0.83	ng/ul	95
22) Benzo(k)fluoranthene	22.81	252	20486	0.74	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.62	276	17914	0.53	ng/ul	98
26) Benzo(g,h,i)perylene	26.30	276	30403	1.12	ng/ul#	92

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

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