

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092216\
 Data File : BM007514.D
 Acq On : 22 Sep 2016 16:52
 Operator : UM/SJ
 Sample : H4830-09DL 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4VH7DL

Quant Time: Sep 22 18:41:55 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 01:26:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	1479	0.40	ng/ul	0.02
2) Naphthalene-d8	10.58	136	5408	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.43	164	3843	0.40	ng/ul	0.02
10) Phenanthrene-d10	17.18	188	8615	0.40	ng/ul	0.02
16) Chrysene-d12	21.35	240	9842	0.40	ng/ul	0.00
20) Perylene-d12	23.62	264	8954	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	1222	0.19	ng/ul	0.03
14) Fluoranthene-d10	19.20	212	4683	0.23	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.63	128	6551	0.52	ng/ul	98
7) Acenaphthylene	14.14	152	6543	0.40	ng/ul	98
8) Acenaphthene	14.49	153	5103	0.40	ng/ul	96
9) Fluorene	15.50	166	7516	0.48	ng/ul	92
11) Pentachlorophenol	16.83	266	3774	1.60	ng/ul	98
12) Phenanthrene	17.21	178	10361	0.44	ng/ul	99
13) Anthracene	17.30	178	6171	0.29	ng/ul	98
17) Pyrene	19.59	202	29297	0.90	ng/ul	97
18) Benzo(a)anthracene	21.33	228	13388	0.46	ng/ul	95
19) Chrysene	21.38	228	18867	0.61	ng/ul	96
22) Benzo(k)fluoranthene	22.98	252	9100	0.30	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.90	276	12629	0.44	ng/ul#	94
26) Benzo(g,h,i)perylene	26.60	276	14205	0.55	ng/ul	99

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

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