

Data Path : Z:\HPCHEM1\BNA_M\Data\bm100615\
 Data File : BM002854.D
 Acq On : 06 Oct 2015 14:42
 Operator : UM/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 06 15:31:19 2015

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM100515.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Oct 06 15:21:07 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.67	152	7972	0.40	ng/ul	-0.02
4) Naphthalene-d8	11.53	136	32738	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.26	164	17290	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.97	188	37842	0.40	ng/ul	-0.02
18) Chrysene-d12	22.06	240	34747	0.40	ng/ul	-0.01
22) Perylene-d12	24.65	264	30324	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.74	96	8361	0.99	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.07	152	18805	0.39	ng/ul	0.00
16) Fluoranthene-d10	19.95	212	37736	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) 1,4-Dioxane	3.79	88	9588	0.99	ng/ul	99
5) Naphthalene	11.58	128	34736	0.40	ng/ul	98
7) 2-Methylnaphthalene	13.14	142	23882	0.40	ng/ul	98
9) Acenaphthylene	14.99	152	26490	0.41	ng/ul	97
10) Acenaphthene	15.31	153	26197	0.40	ng/ul	99
11) Fluorene	16.29	166	29519	0.40	ng/ul	97
13) Pentachlorophenol	17.64	266	959	0.40	ng/ul	97
14) Phenanthrene	18.02	178	46688	0.40	ng/ul	97
15) Anthracene	18.11	178	44553	0.40	ng/ul	97
17) Fluoranthene	19.99	202	52860	0.40	ng/ul	90
19) Pyrene	20.34	202	54695	0.40	ng/ul	97
20) Benzo(a)anthracene	22.04	228	46811	0.40	ng/ul	99
21) Chrysene	22.11	228	45845	0.40	ng/ul	96
23) Benzo(b)fluoranthene	23.85	252	47036	0.42	ng/ul	98
24) Benzo(k)fluoranthene	23.90	252	40390	0.37	ng/ul	95
25) Benzo(a)pyrene	24.53	252	40891	0.40	ng/ul	95
26) Indeno(1,2,3-cd)pyrene	27.37	276	41090	0.40	ng/ul	98
27) Dibenzo(a,h)anthracene	27.36	278	32596	0.40	ng/ul	96
28) Benzo(g,h,i)perylene	28.21	276	35352	0.40	ng/ul	97

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

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