

Data Path : Z:\HPCHEM1\BNA M\Data\bm100515\
 Data File : BM002845.D
 Acq On : 05 Oct 2015 14:51
 Operator : UM/IZ
 Sample : SSTD0.105
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 05 18:43:19 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 : Mon Oct 05 18:40:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.67	152	7467	0.40	ng/ul	-0.02
4) Naphthalene-d8	11.53	136	29071	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.26	164	14768	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.98	188	32348	0.40	ng/ul	0.00
18) Chrysene-d12	22.07	240	28209	0.40	ng/ul	0.00
22) Perylene-d12	24.65	264	23657	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.74	96	1961	0.10	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.07	152	4908	0.11	ng/ul	0.00
16) Fluoranthene-d10	19.96	212	8162	0.10	ng/ul	0.01
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.79	88	2459	0.11	ng/ul	96
5) Naphthalene	11.58	128	7969	0.10	ng/ul	99
7) 2-Methylnaphthalene	13.14	142	5502	0.10	ng/ul	98
9) Acenaphthylene	14.99	152	5600	0.10	ng/ul	96
10) Acenaphthene	15.31	153	5696	0.10	ng/ul	99
11) Fluorene	16.31	166	6389	0.10	ng/ul	82
13) Pentachlorophenol	17.66	266	238	0.15	ng/ul	96
14) Phenanthrene	18.02	178	10027	0.10	ng/ul	99
15) Anthracene	18.11	178	9518	0.10	ng/ul	99
17) Fluoranthene	19.99	202	11410	0.10	ng/ul	93
19) Pyrene	20.34	202	11654	0.10	ng/ul	99
20) Benzo(a)anthracene	22.11	228	9360	0.10	ng/ul	95
21) Chrysene	22.11	228	9454	0.10	ng/ul	98
23) Benzo(b)fluoranthene	23.85	252	9471	0.11	ng/ul	80
24) Benzo(k)fluoranthene	23.91	252	8787	0.10	ng/ul#	79
25) Benzo(a)pyrene	24.53	252	8140	0.10	ng/ul#	77
26) Indeno(1,2,3-cd)pyrene	27.38	276	8458	0.10	ng/ul	97
27) Dibenzo(a,h)anthracene	27.37	278	6747	0.11	ng/ul#	87
28) Benzo(g,h,i)perylene	28.23	276	7352	0.11	ng/ul#	89

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

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