

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004775.D
 Acq On : 25 Mar 2016 13:36
 Operator : UM/SJ
 Sample : SSTD0.839
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 25 15:06:35 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 15:03:08 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1421	5.00	ng/ul	0.00
4) Naphthalene-d8	10.62	136	6245	5.00	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	3808	5.00	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	8545	5.00	ng/ul	0.00
18) Chrysene-d12	21.40	240	8188	5.00	ng/ul	0.00
22) Perylene-d12	23.69	264	6812	5.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	2335	3.39	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	6245	0.88	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	14746	0.86	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.35	88	2773	3.30	ng/ul#	21
5) Naphthalene	10.67	128	10800	0.73	ng/ul	97
7) 2-Methylnaphthalene	12.28	142	7986	0.85	ng/ul	100
9) Acenaphthylene	14.18	152	12859	0.92	ng/ul	95
10) Acenaphthene	14.52	153	9197	0.73	ng/ul	89
11) Fluorene	15.51	166	10916	0.82	ng/ul	90
13) Pentachlorophenol	16.85	266	1518	1.32	ng/ul	96
14) Phenanthrene	17.24	178	17664	0.75	ng/ul	95
15) Anthracene	17.33	178	16738	0.87	ng/ul	95
17) Fluoranthene	19.27	202	20086	0.83	ng/ul	98
19) Pyrene	19.63	202	20773	0.73	ng/ul	98
20) Benzo(a)anthracene	21.38	228	18373	0.95	ng/ul	96
21) Chrysene	21.43	228	18002	0.68	ng/ul	95
23) Benzo(b)fluoranthene	22.99	252	17009	0.70	ng/ul	98
24) Benzo(k)fluoranthene	23.03	252	16397	0.66	ng/ul#	93
25) Benzo(a)pyrene	23.59	252	15570	0.76	ng/ul	97
26) Indeno(1,2,3-cd)pyrene	26.01	276	13893	0.62	ng/ul#	84
27) Dibenzo(a,h)anthracene	26.03	278	10917	0.62	ng/ul#	94
28) Benzo(g,h,i)perylene	26.72	276	11535	0.54	ng/ul	94

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

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