

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004776.D
 Acq On : 25 Mar 2016 14:13
 Operator : UM/SJ
 Sample : SSTD1.640
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 25 15:11:49 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	1227	5.00	ng/ul	0.00
4) Naphthalene-d8	10.62	136	5611	5.00	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	3421	5.00	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	7585	5.00	ng/ul	0.00
18) Chrysene-d12	21.40	240	7547	5.00	ng/ul	0.00
22) Perylene-d12	23.69	264	6874	5.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	4200	7.06	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	11752	1.84	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	28148	1.84	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.35	88	4945	6.82	ng/ul#	22
5) Naphthalene	10.67	128	20460	1.55	ng/ul	99
7) 2-Methylnaphthalene	12.28	142	15173	1.80	ng/ul	99
9) Acenaphthylene	14.18	152	25062	1.99	ng/ul	95
10) Acenaphthene	14.52	153	17450	1.55	ng/ul	89
11) Fluorene	15.51	166	20847	1.75	ng/ul	91
13) Pentachlorophenol	16.85	266	3310	3.25	ng/ul	96
14) Phenanthrene	17.24	178	33405	1.60	ng/ul	95
15) Anthracene	17.33	178	32401	1.90	ng/ul	96
17) Fluoranthene	19.27	202	38362	1.78	ng/ul	98
19) Pyrene	19.63	202	39945	1.52	ng/ul	97
20) Benzo(a)anthracene	21.38	228	36000	2.01	ng/ul	96
21) Chrysene	21.43	228	35127	1.43	ng/ul	97
23) Benzo(b)fluoranthene	22.99	252	35587	1.45	ng/ul	100
24) Benzo(k)fluoranthene	23.03	252	33790	1.36	ng/ul	93
25) Benzo(a)pyrene	23.59	252	33443	1.62	ng/ul	98
26) Indeno(1,2,3-cd)pyrene	26.01	276	32303	1.42	ng/ul#	84
27) Dibenzo(a,h)anthracene	26.03	278	25466	1.44	ng/ul	95
28) Benzo(g,h,i)perylene	26.72	276	26371	1.22	ng/ul	96

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

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