

Data Path : Z:\HPCHEM1\BNA_M\Data\BM012916\
 Data File : BM004145.D
 Acq On : 29 Jan 2016 13:34
 Operator : SJ/IZ
 Sample : SSTD3.253
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.253

Quant Time: Jan 29 14:08:21 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM012916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:08:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1512	0.40	ng/ul	-0.02
4) Naphthalene-d8	10.45	136	6895	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.31	164	4441	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.16	188	544413	0.40	ng/ul	0.09
18) Chrysene-d12	21.28	240	13225	0.40	ng/ul	0.00
22) Perylene-d12	23.37	264	696863	0.40	ng/ul	-0.15
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.12	96	9561	14.79	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.05	152	26362	3.19	ng/ul	-0.01
16) Fluoranthene-d10	19.11	212	79463	0.07	ng/ul	-0.01
Target Compounds				Qvalue		
3) 1,4-Dioxane	3.15	88	11727	14.25	ng/ul#	12
5) Naphthalene	10.50	128	47884	3.08	ng/ul	100
7) 2-Methylnaphthalene	12.12	142	35041	3.36	ng/ul	99
9) Acenaphthylene	14.02	152	56854	3.35	ng/ul	98
10) Acenaphthene	14.37	153	41165	3.01	ng/ul#	76
11) Fluorene	15.38	166	51847	3.42	ng/ul	98
13) Pentachlorophenol	16.73	266	4269	0.05	ng/ul	95
14) Phenanthrene	17.19	178	88707	0.06	ng/ul	93
15) Anthracene	17.19	178	89055	0.07	ng/ul	97
17) Fluoranthene	19.13	202	109268	0.07	ng/ul	93
19) Pyrene	19.51	202	119839	3.36	ng/ul	94
20) Benzo(a)anthracene	21.26	228	113505	3.52	ng/ul	97
21) Chrysene	21.31	228	116241	3.14	ng/ul	96
23) Benzo(b)fluoranthene	22.84	252	121974	0.05	ng/ul	95
24) Benzo(k)fluoranthene	22.84	252	121974	0.06	ng/ul	94
25) Benzo(a)pyrene	23.41	252	118981	0.06	ng/ul#	94
26) Indeno(1,2,3-cd)pyrene	25.76	276	91452	0.04	ng/ul#	63
27) Dibenzo(a,h)anthracene	25.77	278	70249	0.04	ng/ul	97

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

