

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004772.D
 Acq On : 25 Mar 2016 11:48
 Operator : UM/SJ
 Sample : SSTD0.136
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.136

Quant Time: Apr 05 18:34:30 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 18:15:36 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1474	0.40	ng/ul	0.00
4) Naphthalene-d8	10.62	136	6286	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	3882	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	8812	0.40	ng/ul	0.00
18) Chrysene-d12	21.40	240	8470	0.40	ng/ul	0.00
22) Perylene-d12	23.69	264	7467	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	361	0.51	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	962	0.13	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	2198	0.12	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.35	88	451	0.52	ng/ul#	17
5) Naphthalene	10.67	128	1679	0.11	ng/ul#	86
7) 2-Methylnaphthalene	12.28	142	1124	0.12	ng/ul	99
9) Acenaphthylene	14.18	152	1854	0.13	ng/ul#	95
10) Acenaphthene	14.52	153	1301	0.10	ng/ul#	85
11) Fluorene	15.51	166	1577	0.12	ng/ul#	87
14) Phenanthrene	17.24	178	2609	0.11	ng/ul#	94
15) Anthracene	17.33	178	2452	0.12	ng/ul#	92
17) Fluoranthene	19.27	202	2962	0.12	ng/ul	89
19) Pyrene	19.63	202	3142	0.11	ng/ul	95
20) Benzo(a)anthracene	21.38	228	2796	0.14	ng/ul#	90
21) Chrysene	21.43	228	2616	0.09	ng/ul#	89
23) Benzo(b)fluoranthene	23.00	252	2711	0.10	ng/ul#	68
24) Benzo(k)fluoranthene	23.04	252	2401	0.09	ng/ul#	72
25) Benzo(a)pyrene	23.59	252	2395	0.11	ng/ul#	62
26) Indeno(1,2,3-cd)pyrene	26.02	276	2187	0.09	ng/ul#	85
27) Dibenzo(a,h)anthracene	26.04	278	1721	0.09	ng/ul#	61
28) Benzo(a,h,i)perylene	26.73	276	1830	0.08	ng/ul#	70

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

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