

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030915\
 Data File : BM000666.D
 Acq On : 09 Mar 2015 18:21
 Operator : TP/IZ
 Sample : SSTD1.651
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.651

Quant Time: Mar 10 02:48:16 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM030915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 02:25:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4982	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	17204	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4556	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16864	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	14061	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	12004	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.03	152	33558	1.47	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	54909	1.52	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	68516	1.56	ng/ul	98
5) 2-Methylnaphthalene	12.11	142	44876	1.47	ng/ul	99
7) Acenaphthylene	14.03	152	59329	2.77	ng/ul	98
8) Acenaphthene	14.36	153	45032	2.74	ng/ul	98
9) Fluorene	15.36	166	51963	2.59	ng/ul	99
11) Pentachlorophenol	16.71	266	2961	0.71	ng/ul	94
12) Phenanthrene	17.10	178	77970	1.61	ng/ul	99
13) Anthracene	17.17	178	71863m	1.59	ng/ul	
15) Fluoranthene	19.11	202	84880	1.55	ng/ul	99
17) Pyrene	19.48	202	86383	1.38	ng/ul	99
18) Benzo(a)anthracene	21.23	228	67525	1.53	ng/ul	99
19) Chrysene	21.28	228	75928	1.56	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	70665	1.57	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	69777	1.56	ng/ul	94
23) Benzo(a)pyrene	23.35	252	64976	1.59	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.67	276	73765	1.85	ng/ul	99
25) Dibenzo(a,h)anthracene	25.68	278	59852	1.88	ng/ul	94
26) Benzo(g,h,i)perylene	26.33	276	64556	1.85	ng/ul	97

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

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