

Data Path : S:\HPCHEM1\BNA_M\DATA\BM061115\
 Data File : BM001636.D
 Acq On : 10 Jun 2015 21:33
 Operator : TP/IZ
 Sample : SSTD0.185
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.185

Quant Time: Jun 11 07:48:58 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 07:43:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1694	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	5866	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.34	164	2542	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.09	188	5009	0.40	ng/ul	0.00
18) Chrysene-d12	21.27	240	3989	0.40	ng/ul	0.00
22) Perylene-d12	23.50	264	3555	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.08	152	702	0.10	ng/ul	0.00
16) Fluoranthene-d10	19.12	212	1061	0.12	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	10.53	128	1573	0.11	ng/ul#	92
7) 2-Methylnaphthalene	12.15	142	966	0.10	ng/ul	100
9) Acenaphthylene	14.05	152	1226	0.10	ng/ul#	91
10) Acenaphthene	14.39	153	943	0.10	ng/ul	97
11) Fluorene	15.40	166	986	0.09	ng/ul#	80
14) Phenanthrene	17.12	178	1462	0.10	ng/ul#	91
15) Anthracene	17.22	178	1420	0.11	ng/ul#	89
17) Fluoranthene	19.14	202	1651	0.12	ng/ul	92
19) Pyrene	19.50	202	1644	0.10	ng/ul	93
20) Benzo(a)anthracene	21.25	228	1334	0.12	ng/ul	93
21) Chrysene	21.31	228	1496	0.11	ng/ul	96
23) Benzo(b)fluoranthene	22.83	252	1490	0.12	ng/ul	83
24) Benzo(k)fluoranthene	22.87	252	1662	0.13	ng/ul#	83
25) Benzo(a)pyrene	23.40	252	1438	0.12	ng/ul#	77
26) Indeno(1,2,3-cd)pyrene	25.77	276	1820	0.14	ng/ul	98
27) Dibenzo(a,h)anthracene	25.78	278	1506	0.15	ng/ul#	85
28) Benzo(g,h,i)perylene	26.45	276	1774	0.16	ng/ul	93

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

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