

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030915\
 Data File : BM000663.D
 Acq On : 09 Mar 2015 16:34
 Operator : TP/IZ
 Sample : SSTD0.248
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.248

Quant Time: Mar 10 02:36:33 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	4790	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	16270	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4683	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15575	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	12481	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	11429	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	4112	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6478	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	8540	0.21	ng/ul	99
5) 2-Methylnaphthalene	12.12	142	5374	0.19	ng/ul	100
7) Acenaphthylene	14.03	152	6739	0.31	ng/ul	100
8) Acenaphthene	14.36	153	5019	0.30	ng/ul	99
9) Fluorene	15.36	166	6073	0.29	ng/ul	98
11) Pentachlorophenol	16.72	266	177	0.05	ng/ul	90
12) Phenanthrene	17.10	178	9058	0.20	ng/ul	99
13) Anthracene	17.19	178	7943	0.19	ng/ul	99
15) Fluoranthene	19.11	202	9686	0.19	ng/ul	99
17) Pyrene	19.48	202	9828	0.18	ng/ul	98
18) Benzo(a)anthracene	21.23	228	7499	0.19	ng/ul	100
19) Chrysene	21.28	228	9249	0.21	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	8146	0.19	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	8827	0.21	ng/ul	95
23) Benzo(a)pyrene	23.35	252	7803	0.20	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.67	276	8992	0.24	ng/ul	99
25) Dibenzo(a,h)anthracene	25.68	278	7232	0.24	ng/ul	98
26) Benzo(g,h,i)perylene	26.35	276	8162	0.25	ng/ul	99

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

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