

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042715\
 Data File : BM001113.D
 Acq On : 27 Apr 2015 11:17
 Operator : TP/IZ
 Sample : SSTD0.244
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.244

Quant Time: Apr 27 14:12:00 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM042715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 27 13:50:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	1086	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3756	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	1893	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	4234	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3944	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	3821	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.93	152	931	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1938	0.21	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.37	128	1886	0.20	ng/ul	94
5) 2-Methylnaphthalene	12.00	142	1216	0.19	ng/ul	98
7) Acenaphthylene	13.91	152	1697	0.20	ng/ul	96
8) Acenaphthene	14.27	153	1312	0.20	ng/ul	98
9) Fluorene	15.26	166	1483	0.19	ng/ul	97
11) Pentachlorophenol	16.63	266	155	0.16	ng/ul	94
12) Phenanthrene	17.00	178	2408	0.20	ng/ul	97
13) Anthracene	17.09	178	2184	0.20	ng/ul	97
15) Fluoranthene	19.02	202	2714	0.20	ng/ul	97
17) Pyrene	19.38	202	2860	0.19	ng/ul	97
18) Benzo(a)anthracene	21.14	228	2440	0.20	ng/ul	98
19) Chrysene	21.19	228	2762	0.20	ng/ul	98
21) Benzo(b)fluoranthene	22.67	252	2637	0.19	ng/ul	93
22) Benzo(k)fluoranthene	22.71	252	2779	0.20	ng/ul	94
23) Benzo(a)pyrene	23.22	252	2649	0.20	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.46	276	2986	0.20	ng/ul	99
25) Dibenzo(a,h)anthracene	25.49	278	2392	0.20	ng/ul	93
26) Benzo(g,h,i)perylene	26.12	276	2619	0.20	ng/ul	95

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

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