

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041715\
 Data File : BM001023.D
 Acq On : 16 Apr 2015 18:57
 Operator : TP/IZ
 Sample : SSTD0.228
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.228

Quant Time: Apr 17 06:36:05 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM041715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 06:23:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1382	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	4616	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2872	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	4697	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	3934	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	3688	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.97	152	1130	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	1993	0.22	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	2356	0.22	ng/ul	96
5) 2-Methylnaphthalene	12.05	142	1503	0.22	ng/ul	99
7) Acenaphthylene	13.96	152	1922	0.20	ng/ul	96
8) Acenaphthene	14.30	153	1537	0.20	ng/ul	99
9) Fluorene	15.31	166	1696	0.19	ng/ul#	79
11) Pentachlorophenol	16.67	266	53	0.10	ng/ul	92
12) Phenanthrene	17.02	178	2663	0.21	ng/ul	97
13) Anthracene	17.12	178	2329	0.21	ng/ul	94
15) Fluoranthene	19.05	202	2919	0.21	ng/ul	97
17) Pyrene	19.42	202	2979	0.21	ng/ul#	92
18) Benzo(a)anthracene	21.18	228	2297	0.21	ng/ul	98
19) Chrysene	21.23	228	2906	0.23	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	2530	0.21	ng/ul	92
22) Benzo(k)fluoranthene	22.75	252	2790	0.22	ng/ul	93
23) Benzo(a)pyrene	23.26	252	2544	0.22	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.53	276	3250	0.24	ng/ul	99
25) Dibenzo(a,h)anthracene	25.54	278	2681	0.25	ng/ul	95
26) Benzo(g,h,i)perylene	26.19	276	2895	0.24	ng/ul	97

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

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