

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051415\
 Data File : BM001280.D
 Acq On : 13 May 2015 18:24
 Operator : TP/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.472

Quant Time: May 14 01:34:56 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM051115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 14 01:20:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	1064	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	3717	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	1891	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	3858	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	3033	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	2805	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.83	152	1900	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.91	212	3234	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.27	128	3786	0.40	ng/ul	99
5) 2-Methylnaphthalene	11.90	142	2461	0.40	ng/ul	100
7) Acenaphthylene	13.82	152	3339	0.40	ng/ul	97
8) Acenaphthene	14.17	153	2633	0.40	ng/ul	91
9) Fluorene	15.18	166	2911	0.40	ng/ul	91
11) Pentachlorophenol	16.53	266	178	0.21	ng/ul	94
12) Phenanthrene	16.91	178	4374	0.40	ng/ul	97
13) Anthracene	17.00	178	3958	0.39	ng/ul	99
15) Fluoranthene	18.94	202	4751	0.39	ng/ul	99
17) Pyrene	19.30	202	4846	0.41	ng/ul	99
18) Benzo(a)anthracene	21.06	228	3775	0.41	ng/ul	98
19) Chrysene	21.11	228	4308	0.41	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	4200	0.42	ng/ul	99
22) Benzo(k)fluoranthene	22.60	252	4147	0.40	ng/ul	99
23) Benzo(a)pyrene	23.09	252	3885	0.40	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.27	276	4563	0.43	ng/ul	99
25) Dibenzo(a,h)anthracene	25.28	278	3614	0.42	ng/ul	97
26) Benzo(g,h,i)perylene	25.90	276	4114	0.43	ng/ul	99

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

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