

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051415\
 Data File : BM001276.D
 Acq On : 13 May 2015 15:28
 Operator : TP/IZ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.471

Quant Time: May 14 01:27:49 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 : Tue May 12 03:03:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	1224	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	4335	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.12	164	2323	0.40	ng/ul	-0.02
10) Phenanthrene-d10	16.86	188	4629	0.40	ng/ul	-0.02
16) Chrysene-d12	21.08	240	3577	0.40	ng/ul	-0.01
20) Perylene-d12	23.19	264	3259	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.83	152	2272	0.41	ng/ul	-0.01
14) Fluoranthene-d10	18.91	212	3823	0.38	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.27	128	4486	0.41	ng/ul	99
5) 2-Methylnaphthalene	11.90	142	2965	0.41	ng/ul	100
7) Acenaphthylene	13.82	152	4135	0.40	ng/ul	96
8) Acenaphthene	14.17	153	3191	0.39	ng/ul	89
9) Fluorene	15.18	166	3556	0.39	ng/ul	90
11) Pentachlorophenol	16.53	266	305	0.30	ng/ul	91
12) Phenanthrene	16.91	178	5218	0.40	ng/ul	96
13) Anthracene	17.00	178	4849	0.40	ng/ul	98
15) Fluoranthene	18.94	202	5645	0.39	ng/ul	97
17) Pyrene	19.30	202	5759	0.41	ng/ul	96
18) Benzo(a)anthracene	21.06	228	4568	0.42	ng/ul	95
19) Chrysene	21.11	228	5121	0.41	ng/ul	95
21) Benzo(b)fluoranthene	22.56	252	4583	0.39	ng/ul	98
22) Benzo(k)fluoranthene	22.60	252	5044	0.42	ng/ul	98
23) Benzo(a)pyrene	23.10	252	4553	0.40	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.28	276	5213	0.42	ng/ul	99
25) Dibenzo(a,h)anthracene	25.29	278	4167	0.42	ng/ul	98
26) Benzo(g,h,i)perylene	25.91	276	4606	0.42	ng/ul	99

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

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