

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041715\
 Data File : BM001039.D
 Acq On : 17 Apr 2015 19:45
 Operator : TP/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.439

Quant Time: Apr 18 00:14:50 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 15:43:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1445	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	5165	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2661	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	5702	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	4830	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	4373	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.96	152	2620	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	4865	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.41	128	5243	0.39	ng/ul	97
5) 2-Methylnaphthalene	12.03	142	3462	0.39	ng/ul	100
7) Acenaphthylene	13.94	152	4503	0.38	ng/ul	99
8) Acenaphthene	14.30	153	3632	0.39	ng/ul	94
9) Fluorene	15.30	166	4206	0.39	ng/ul	86
11) Pentachlorophenol	16.64	266	374	0.28	ng/ul	91
12) Phenanthrene	17.03	178	6465	0.39	ng/ul	94
13) Anthracene	17.12	178	5860	0.39	ng/ul	94
15) Fluoranthene	19.06	202	7117	0.39	ng/ul	90
17) Pyrene	19.42	202	7298	0.40	ng/ul	96
18) Benzo(a)anthracene	21.17	228	5788	0.38	ng/ul	96
19) Chrysene	21.22	228	6721	0.40	ng/ul	97
21) Benzo(b)fluoranthene	22.71	252	6548	0.41	ng/ul	98
22) Benzo(k)fluoranthene	22.75	252	6168	0.38	ng/ul	98
23) Benzo(a)pyrene	23.27	252	5859	0.39	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.53	276	6679	0.40	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	5435	0.40	ng/ul	99
26) Benzo(g,h,i)perylene	26.18	276	5930	0.40	ng/ul	97

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

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