

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM102516\
 Data File : BM007731.D
 Acq On : 25 Oct 2016 09:12
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.461

Quant Time: Oct 25 11:24:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 25 11:17:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	1060	0.40	ng/ul	0.00
2) Naphthalene-d8	10.55	136	3373	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	1539	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.14	188	3349	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	3191	0.40	ng/ul	0.00
20) Perylene-d12	23.59	264	2957	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.15	152	1629	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	3412	0.45	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.60	128	3521	0.41	ng/ul	99
5) 2-Methylnaphthalene	12.23	142	2241	0.41	ng/ul	100
7) Acenaphthylene	14.11	152	3019	0.42	ng/ul	99
8) Acenaphthene	14.45	153	2329	0.42	ng/ul	100
9) Fluorene	15.46	166	2556	0.42	ng/ul	99
11) Pentachlorophenol	16.82	266	316	0.40	ng/ul	99
12) Phenanthrene	17.18	178	3963	0.42	ng/ul	99
13) Anthracene	17.28	178	3639	0.43	ng/ul	98
15) Fluoranthene	19.21	202	4954	0.44	ng/ul	99
17) Pyrene	19.57	202	5022	0.46	ng/ul	98
18) Benzo(a)anthracene	21.31	228	3982	0.44	ng/ul	99
19) Chrysene	21.37	228	4738	0.46	ng/ul	99
21) Benzo(b)fluoranthene	22.90	252	4905	0.48	ng/ul	97
22) Benzo(k)fluoranthene	22.95	252	4386	0.42	ng/ul	100
23) Benzo(a)pyrene	23.49	252	4193	0.44	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.86	276	5206	0.46	ng/ul	100
25) Dibenzo(a,h)anthracene	25.88	278	4152	0.45	ng/ul	99
26) Benzo(g,h,i)perylene	26.56	276	4642	0.45	ng/ul	100

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

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