

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006212.D
 Acq On : 09 Jun 2019 04:15
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.404

Quant Time: Jun 09 05:13:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	7610	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	34577	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	17408	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	33386	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	24280	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	27206	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	20191	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	34201	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	37519	0.393	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	25517	0.384	ng/ul	99
7) Acenaphthylene	14.01	152	37104	0.383	ng/ul	99
8) Acenaphthene	14.36	153	26796	0.399	ng/ul	100
9) Fluorene	15.35	166	28240	0.360	ng/ul	97
11) Pentachlorophenol	16.71	266	3391	0.400	ng/ul	99
12) Phenanthrene	17.09	178	40857	0.381	ng/ul	100
13) Anthracene	17.18	178	37206	0.364	ng/ul	100
15) Fluoranthene	19.12	202	44849	0.379	ng/ul	97
17) Pyrene	19.48	202	46516	0.364	ng/ul	98
18) Benzo(a)anthracene	21.24	228	39019	0.363	ng/ul	98
19) Chrysene	21.29	228	43496	0.431	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	41917	0.370	ng/ul	97
22) Benzo(k)fluoranthene	22.86	252	43646	0.395	ng/ul	95
23) Benzo(a)pyrene	23.39	252	42356	0.401	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.71	276	48744	0.426	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.73	278	38690	0.423	ng/ul	93
26) Benzo(g,h,i)perylene	26.40	276	41189	0.430	ng/ul	94

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

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