

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN071019\
 Data File : BN006773.D
 Acq On : 10 Jul 2019 17:05
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Jul 11 01:05:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 11 01:03:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2176	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	8378	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4441	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9481	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	8194	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	9326	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4842	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	10651	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.42	128	9599	0.399	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	5886	0.362	ng/ul	99
7) Acenaphthylene	12.97	152	8384	0.366	ng/ul	99
8) Acenaphthene	14.31	153	6908	0.391	ng/ul	99
9) Fluorene	15.31	166	7134	0.368	ng/ul	99
11) Pentachlorophenol	16.71	266	520	0.368	ng/ul#	84
12) Phenanthrene	17.06	178	11606	0.381	ng/ul	99
13) Anthracene	17.15	178	8970	0.340	ng/ul	99
15) Fluoranthene	19.09	202	13882	0.368	ng/ul	97
17) Pyrene	19.46	202	14791	0.395	ng/ul	97
18) Benzo(a)anthracene	21.23	228	11698	0.348	ng/ul	99
19) Chrysene	21.28	228	15726	0.419	ng/ul	100
21) Benzo(b)fluoranthene	22.81	252	14351	0.375	ng/ul	97
22) Benzo(k)fluoranthene	22.86	252	15790	0.395	ng/ul	98
23) Benzo(a)pyrene	23.39	252	14359	0.384	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.71	276	16755	0.442	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.71	278	13090	0.434	ng/ul	97
26) Benzo(g,h,i)perylene	26.39	276	14736	0.465	ng/ul	97

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

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