

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092218\
 Data File : BN002829.D
 Acq On : 21 Sep 2018 21:39
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.473

Quant Time: Sep 21 23:49:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 21 23:48:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1065	0.40	ng/ul	0.01
2) Naphthalene-d8	10.44	136	4723	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	2461	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	5310	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4812	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4550	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	2821	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	5907	0.43	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.49	128	4895	0.412	ng/ul#	96
5) 2-Methylnaphthalene	12.12	142	3241	0.421	ng/ul	99
7) Acenaphthylene	14.01	152	4250	0.409	ng/ul	99
8) Acenaphthene	14.36	153	3479	0.410	ng/ul	98
9) Fluorene	15.36	166	3820	0.422	ng/ul	93
11) Pentachlorophenol	16.74	266	255	0.426	ng/ul	95
12) Phenanthrene	17.09	178	6004	0.410	ng/ul#	89
13) Anthracene	17.20	178	5320	0.418	ng/ul#	93
15) Fluoranthene	19.12	202	6720	0.426	ng/ul	96
17) Pyrene	19.48	202	6840	0.436	ng/ul#	96
18) Benzo(a)anthracene	21.24	228	5790	0.435	ng/ul#	86
19) Chrysene	21.29	228	6649	0.429	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	6288	0.418	ng/ul	90
22) Benzo(k)fluoranthene	22.85	252	6616	0.418	ng/ul	92
23) Benzo(a)pyrene	23.38	252	6137	0.426	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.69	276	6981	0.402	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.70	278	5644	0.406	ng/ul	95
26) Benzo(g,h,i)perylene	26.37	276	6011	0.401	ng/ul	99

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

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