

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120118\
 Data File : BN003675.D
 Acq On : 02 Dec 2018 00:04
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.479

Quant Time: Dec 03 02:03:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 03 02:00:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4964	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	19933	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	10458	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	23982	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	23070	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	20716	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	12050	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	27176	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.69	128	22609	0.404	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	15876	0.406	ng/ul	99
7) Acenaphthylene	14.20	152	19781	0.420	ng/ul	100
8) Acenaphthene	14.54	153	16804	0.407	ng/ul	99
9) Fluorene	15.52	166	19069	0.396	ng/ul	99
11) Pentachlorophenol	16.87	266	1967	0.455	ng/ul	99
12) Phenanthrene	17.26	178	31992	0.406	ng/ul	99
13) Anthracene	17.35	178	27004	0.417	ng/ul	100
15) Fluoranthene	19.27	202	36830	0.396	ng/ul	98
17) Pyrene	19.63	202	35898	0.434	ng/ul	97
18) Benzo(a)anthracene	21.38	228	30667	0.421	ng/ul	99
19) Chrysene	21.43	228	33679	0.404	ng/ul	99
21) Benzo(b)fluoranthene	23.00	252	34718	0.405	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	33761	0.404	ng/ul	99
23) Benzo(a)pyrene	23.61	252	30268	0.408	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	32383	0.381	ng/ul	99
25) Dibenzo(a,h)anthracene	26.07	278	25366	0.370	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	27601	0.373	ng/ul	99

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

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