

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121118\
 Data File : BN003815.D
 Acq On : 11 Dec 2018 15:25
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.498

Quant Time: Dec 11 16:00:47 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 : Tue Dec 11 15:12:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4500	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	18077	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8849	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	17518	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12433	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	12202	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	10142	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	16479	0.32	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.68	128	19424	0.383	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	13283	0.375	ng/ul	99
7) Acenaphthylene	14.19	152	17257	0.433	ng/ul	100
8) Acenaphthene	14.53	153	13417	0.384	ng/ul	99
9) Fluorene	15.52	166	14487	0.356	ng/ul	100
11) Pentachlorophenol	16.86	266	1165	0.369	ng/ul	98
12) Phenanthrene	17.26	178	21687	0.377	ng/ul	99
13) Anthracene	17.35	178	19530	0.413	ng/ul	99
15) Fluoranthene	19.27	202	22024	0.324	ng/ul	97
17) Pyrene	19.63	202	21244	0.476	ng/ul	96
18) Benzo(a)anthracene	21.38	228	17277	0.440	ng/ul	98
19) Chrysene	21.44	228	17122	0.381	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	17383	0.344	ng/ul	94
22) Benzo(k)fluoranthene	23.06	252	17602	0.358	ng/ul#	95
23) Benzo(a)pyrene	23.61	252	16938	0.388	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.07	276	20771	0.414	ng/ul	98
25) Dibenzo(a,h)anthracene	26.08	278	16495	0.409	ng/ul	96
26) Benzo(g,h,i)perylene	26.80	276	17722	0.407	ng/ul	96

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

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