

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003612.D
 Acq On : 20 Nov 2018 01:41
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.469

Quant Time: Nov 20 03:22:12 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 Nov 20 01:46:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5285	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20812	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11378	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	27583	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	33251	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	28937	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	12890	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	32961	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.69	128	23565	0.403	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	16557	0.406	ng/ul	100
7) Acenaphthylene	14.20	152	22392	0.437	ng/ul	100
8) Acenaphthene	14.54	153	17922	0.399	ng/ul	99
9) Fluorene	15.53	166	20921	0.400	ng/ul	99
11) Pentachlorophenol	16.87	266	2162	0.435	ng/ul	100
12) Phenanthrene	17.26	178	36403	0.402	ng/ul	99
13) Anthracene	17.35	178	30804	0.414	ng/ul	100
15) Fluoranthene	19.28	202	43152	0.404	ng/ul	100
17) Pyrene	19.64	202	42683	0.358	ng/ul	99
18) Benzo(a)anthracene	21.38	228	44004	0.419	ng/ul	99
19) Chrysene	21.43	228	46311	0.385	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	46773	0.391	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	45932	0.394	ng/ul	100
23) Benzo(a)pyrene	23.61	252	41821	0.403	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	45750	0.385	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.08	278	36825	0.385	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	38196	0.370	ng/ul	99

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

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