

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003380.D
 Acq On : 01 Nov 2018 16:03
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.432

Quant Time: Nov 01 17:51:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:18:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	7758	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	30016	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	17596	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	42829	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	43660	0.40	ng/ul	0.00
20) Perylene-d12	23.74	264	38250	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	19686	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	48211	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.70	128	34263	0.407	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	24611	0.406	ng/ul	100
7) Acenaphthylene	14.21	152	30118	0.394	ng/ul	99
8) Acenaphthene	14.55	153	27266	0.399	ng/ul	99
9) Fluorene	15.54	166	31856	0.396	ng/ul	99
11) Pentachlorophenol	16.87	266	3119	0.364	ng/ul	98
12) Phenanthrene	17.27	178	55698	0.397	ng/ul	99
13) Anthracene	17.36	178	45103	0.394	ng/ul	100
15) Fluoranthene	19.29	202	62648	0.397	ng/ul	99
17) Pyrene	19.66	202	62053	0.391	ng/ul	99
18) Benzo(a)anthracene	21.40	228	54545	0.398	ng/ul	100
19) Chrysene	21.46	228	62634	0.403	ng/ul	100
21) Benzo(b)fluoranthene	23.04	252	66027	0.401	ng/ul	99
22) Benzo(k)fluoranthene	23.08	252	62771	0.393	ng/ul	99
23) Benzo(a)pyrene	23.64	252	58185	0.423	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.10	276	59983	0.407	ng/ul	100
25) Dibenzo(a,h)anthracene	26.11	278	48620	0.405	ng/ul	99
26) Benzo(g,h,i)perylene	26.82	276	49411	0.391	ng/ul	100

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

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