

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003138.D
 Acq On : 12 Oct 2018 13:07
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.407

Quant Time: Oct 12 13:49:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1083	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	5171	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3076	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7058	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	4963	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4397	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3533	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8219	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	5684	0.428	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	4029	0.441	ng/ul	100
7) Acenaphthylene	13.99	152	6128	0.435	ng/ul	99
8) Acenaphthene	14.33	153	4709	0.441	ng/ul	99
9) Fluorene	15.34	166	5433	0.444	ng/ul	98
11) Pentachlorophenol	16.74	266	397	0.580	ng/ul	91
12) Phenanthrene	17.07	178	8595	0.437	ng/ul	99
13) Anthracene	17.17	178	8100	0.436	ng/ul	100
15) Fluoranthene	19.09	202	9347	0.405	ng/ul	98
17) Pyrene	19.46	202	9271	0.481	ng/ul	99
18) Benzo(a)anthracene	21.22	228	7003	0.438	ng/ul	100
19) Chrysene	21.27	228	7192	0.443	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	6410	0.405	ng/ul	99
22) Benzo(k)fluoranthene	22.84	252	6874	0.419	ng/ul	99
23) Benzo(a)pyrene	23.36	252	6256	0.441	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.68	276	7417	0.455	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.68	278	5951	0.456	ng/ul	96
26) Benzo(g,h,i)perylene	26.35	276	6290	0.457	ng/ul	99

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

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