

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003592.D
 Acq On : 19 Nov 2018 14:00
 Operator : MJ/SJ
 Sample : SSTD0.864
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.864

Quant Time: Nov 19 14:36:51 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 Nov 19 13:56:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	5372	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20346	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11147	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	26805	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	30262	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	26713	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	25903	0.80	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	64723	0.86	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.69	128	46972	0.823	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	33191	0.809	ng/ul	99
7) Acenaphthylene	14.20	152	41048	0.847	ng/ul	100
8) Acenaphthene	14.54	153	35862	0.829	ng/ul	99
9) Fluorene	15.53	166	42086	0.826	ng/ul	100
11) Pentachlorophenol	16.87	266	3769	0.702	ng/ul	99
12) Phenanthrene	17.26	178	72765	0.828	ng/ul	99
13) Anthracene	17.36	178	60558	0.845	ng/ul	100
15) Fluoranthene	19.28	202	87120	0.882	ng/ul	99
17) Pyrene	19.64	202	84431	0.767	ng/ul	100
18) Benzo(a)anthracene	21.38	228	77728	0.819	ng/ul	99
19) Chrysene	21.43	228	88442	0.820	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	90928	0.790	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	89888	0.806	ng/ul	98
23) Benzo(a)pyrene	23.61	252	79500	0.827	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.06	276	91730	0.892	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	74216	0.885	ng/ul	98
26) Benzo(g,h,i)perylene	26.78	276	78804	0.893	ng/ul	98

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

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