

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111918\
 Data File : BN003597.D
 Acq On : 19 Nov 2018 16:56
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.468

Quant Time: Nov 19 17:36:08 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 15:40:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	4851	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	19206	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11068	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	27534	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	24482	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	21601	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	12573	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	26804	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.69	128	21042	0.390	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	15978	0.424	ng/ul	99
7) Acenaphthylene	14.20	152	19289	0.387	ng/ul	100
8) Acenaphthene	14.54	153	17321	0.397	ng/ul	99
9) Fluorene	15.53	166	20165	0.396	ng/ul	100
11) Pentachlorophenol	16.87	266	1621	0.327	ng/ul	99
12) Phenanthrene	17.26	178	35476	0.393	ng/ul	99
13) Anthracene	17.36	178	28335	0.382	ng/ul	100
15) Fluoranthene	19.28	202	35562	0.333	ng/ul	100
17) Pyrene	19.64	202	34685	0.395	ng/ul	99
18) Benzo(a)anthracene	21.38	228	29713	0.384	ng/ul	99
19) Chrysene	21.43	228	35542	0.402	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	36317	0.406	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	34947	0.401	ng/ul	100
23) Benzo(a)pyrene	23.61	252	30722	0.397	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	34671	0.391	ng/ul	100
25) Dibenzo(a,h)anthracene	26.08	278	27798	0.389	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	29969	0.389	ng/ul	100

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

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