

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111918\
 Data File : BN003589.D
 Acq On : 19 Nov 2018 12:15
 Operator : MJ/SJ
 Sample : SSTD0.161
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.161

Quant Time: Nov 19 14:04:17 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	7.84	152	5069	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	18986	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	10372	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	26910	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	28337	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	27631	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	2849	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	7899	0.10	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.69	128	5332	0.100	ng/ul	97
5) 2-Methylnaphthalene	12.30	142	3637	0.095	ng/ul	99
7) Acenaphthylene	14.20	152	4587	0.102	ng/ul	97
8) Acenaphthene	14.54	153	4060	0.101	ng/ul	100
9) Fluorene	15.53	166	4673	0.099	ng/ul	98
11) Pentachlorophenol	16.87	266	348	0.065	ng/ul	99
12) Phenanthrene	17.26	178	8541	0.097	ng/ul	97
13) Anthracene	17.36	178	6834	0.095	ng/ul	98
15) Fluoranthene	19.28	202	10312	0.104	ng/ul	97
17) Pyrene	19.64	202	10238	0.099	ng/ul	98
18) Benzo(a)anthracene	21.38	228	9041	0.102	ng/ul	99
19) Chrysene	21.43	228	10392	0.103	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	11014	0.093	ng/ul	90
22) Benzo(k)fluoranthene	23.06	252	10677	0.093	ng/ul#	90
23) Benzo(a)pyrene	23.61	252	9399	0.095	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.07	276	10832	0.102	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.08	278	8688	0.100	ng/ul	95
26) Benzo(g,h,i)perylene	26.79	276	9495	0.104	ng/ul	97

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

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