

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN002987.D
 Acq On : 08 Oct 2018 15:16
 Operator : MJ/SJ
 Sample : SSTD0.288
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.288

Quant Time: Oct 08 16:29:01 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 : Mon Oct 08 16:26:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1125	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	5214	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.28	164	3258	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	7751	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6959	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	5259	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1590	0.21	ng/ul	0.02
14) Fluoranthene-d10	19.07	212	4289	0.21	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.47	128	2660	0.203	ng/ul#	92
5) 2-Methylnaphthalene	12.11	142	1785	0.210	ng/ul	99
7) Acenaphthylene	14.00	152	2880	0.209	ng/ul	96
8) Acenaphthene	14.34	153	2168	0.193	ng/ul	97
9) Fluorene	15.35	166	2495	0.208	ng/ul	99
11) Pentachlorophenol	16.74	266	136	0.156	ng/ul#	86
12) Phenanthrene	17.08	178	4113	0.192	ng/ul	98
13) Anthracene	17.17	178	3964	0.213	ng/ul	98
15) Fluoranthene	19.10	202	4899	0.213	ng/ul	98
17) Pyrene	19.47	202	4982	0.220	ng/ul	99
18) Benzo(a)anthracene	21.22	228	4386	0.228	ng/ul	97
19) Chrysene	21.27	228	4439	0.198	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	3767	0.217	ng/ul	89
22) Benzo(k)fluoranthene	22.83	252	3894	0.213	ng/ul#	90
23) Benzo(a)pyrene	23.36	252	3315	0.199	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.67	276	3794	0.189	ng/ul	97
25) Dibenzo(a,h)anthracene	25.67	278	3039	0.189	ng/ul#	88
26) Benzo(g,h,i)perylene	26.35	276	3231	0.187	ng/ul	91

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

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