

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100918\
 Data File : BN002991.D
 Acq On : 08 Oct 2018 17:36
 Operator : MJ/SJ
 Sample : SSTD3.292
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.292

Quant Time: Oct 08 18:06:32 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	7.64	152	1316	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	5657	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	3290	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7646	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7355	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5417	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	29015	3.56	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	76068	3.86	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.44	128	46446	3.265	ng/ul 96
5) 2-Methylnaphthalene	12.07	142	33697	3.654	ng/ul 97
7) Acenaphthylene	13.99	152	52108	3.752	ng/ul 97
8) Acenaphthene	14.33	153	37864	3.340	ng/ul 99
9) Fluorene	15.32	166	44346	3.665	ng/ul 99
11) Pentachlorophenol	16.71	266	3386	3.929	ng/ul 96
12) Phenanthrene	17.06	178	72277	3.424	ng/ul 99
13) Anthracene	17.15	178	71823	3.919	ng/ul 98
15) Fluoranthene	19.09	202	86157	3.789	ng/ul 97
17) Pyrene	19.45	202	86261	3.596	ng/ul 97
18) Benzo(a)anthracene	21.21	228	79937	3.931	ng/ul 98
19) Chrysene	21.26	228	76141	3.212	ng/ul 98
21) Benzo(b)fluoranthene	22.78	252	64957	3.628	ng/ul 90
22) Benzo(k)fluoranthene	22.82	252	66507	3.531	ng/ul# 89
23) Benzo(a)pyrene	23.35	252	58397	3.406	ng/ul# 85
24) Indeno(1,2,3-cd)pyrene	25.65	276	60576	2.931	ng/ul# 99
25) Dibenzo(a,h)anthracene	25.65	278	48725	2.944	ng/ul# 88
26) Benzo(g,h,i)perylene	26.33	276	50408	2.825	ng/ul 91

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

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