

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008175.D
 Acq On : 15 Oct 2019 16:15
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Oct 15 16:48:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1948	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.38	136	10223	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.25	164	6746	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.00	188	16522	0.40	ng/ul	-0.02
16) Chrysene-d12	21.20	240	15575	0.40	ng/ul	-0.01
20) Perylene-d12	23.38	264	12891	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	6408	0.37	ng/ul	-0.02
14) Fluoranthene-d10	19.04	212	19511	0.34	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	10736	0.369	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	7506	0.376	ng/ul	99
7) Acenaphthylene	13.97	152	11803	0.367	ng/ul	100
8) Acenaphthene	14.32	153	10273	0.411	ng/ul	99
9) Fluorene	15.31	166	10535	0.371	ng/ul	99
11) Pentachlorophenol	16.67	266	1275	0.319	ng/ul	99
12) Phenanthrene	17.04	178	17399	0.360	ng/ul	99
13) Anthracene	17.14	178	15348	0.361	ng/ul	100
15) Fluoranthene	19.06	202	22018	0.345	ng/ul	98
17) Pyrene	19.43	202	25422	0.410	ng/ul	99
18) Benzo(a)anthracene	21.18	228	20767	0.367	ng/ul	99
19) Chrysene	21.24	228	24369	0.351	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	18599	0.443	ng/ul	96
22) Benzo(k)fluoranthene	22.78	252	20504	0.431	ng/ul	99
23) Benzo(a)pyrene	23.29	252	18125	0.410	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.56	276	18268	0.351	ng/ul	99
25) Dibenzo(a,h)anthracene	25.57	278	14624	0.356	ng/ul	100
26) Benzo(g,h,i)perylene	26.21	276	16011	0.336	ng/ul	98

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

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