

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006987.D
 Acq On : 07 Aug 2019 14:58
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.427

Quant Time: Aug 07 17:41:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	5693	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	21651	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	13476	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	33563	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	29780	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	30432	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	14662	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	40699	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.28	128	23213	0.373	ng/ul	98
5) 2-Methylnaphthalene	11.91	142	17374	0.371	ng/ul	100
7) Acenaphthylene	13.83	152	28824	0.369	ng/ul	99
8) Acenaphthene	14.18	153	19806	0.371	ng/ul	99
9) Fluorene	15.17	166	23945	0.370	ng/ul	100
11) Pentachlorophenol	16.52	266	2993	0.324	ng/ul#	85
12) Phenanthrene	16.90	178	39984	0.374	ng/ul	100
13) Anthracene	16.99	178	40192	0.367	ng/ul	100
15) Fluoranthene	18.93	202	49139	0.370	ng/ul	99
17) Pyrene	19.29	202	50081	0.377	ng/ul	99
18) Benzo(a)anthracene	21.06	228	48494	0.371	ng/ul	99
19) Chrysene	21.11	228	45485	0.371	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	45220	0.367	ng/ul	99
22) Benzo(k)fluoranthene	22.62	252	45094	0.377	ng/ul	98
23) Benzo(a)pyrene	23.12	252	42662	0.367	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.31	276	45349	0.362	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.33	278	36384	0.364	ng/ul	100
26) Benzo(g,h,i)perylene	25.94	276	38127	0.363	ng/ul	100

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

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