

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080919\
 Data File : BN007042.D
 Acq On : 09 Aug 2019 17:18
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.432

Quant Time: Aug 09 17:55:02 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	7.46	152	10532	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.22	136	38821	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.11	164	23315	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	57115	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	44370	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	43712	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	25706	0.37	ng/ul	-0.01
14) Fluoranthene-d10	18.90	212	64813	0.35	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.27	128	41547	0.372	ng/ul	97
5) 2-Methylnaphthalene	11.90	142	30771	0.367	ng/ul	100
7) Acenaphthylene	13.82	152	47541	0.352	ng/ul	99
8) Acenaphthene	14.17	153	34306	0.371	ng/ul	100
9) Fluorene	15.17	166	40836	0.364	ng/ul	99
11) Pentachlorophenol	16.52	266	5451	0.346	ng/ul#	81
12) Phenanthrene	16.90	178	68615	0.377	ng/ul	100
13) Anthracene	16.99	178	64802	0.348	ng/ul	99
15) Fluoranthene	18.93	202	79463	0.351	ng/ul	98
17) Pyrene	19.29	202	79606	0.402	ng/ul	100
18) Benzo(a)anthracene	21.07	228	71672	0.368	ng/ul	100
19) Chrysene	21.12	228	67778	0.371	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	63255	0.357	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	61696	0.359	ng/ul	98
23) Benzo(a)pyrene	23.12	252	59614	0.357	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.31	276	66129	0.368	ng/ul	99
25) Dibenzo(a,h)anthracene	25.33	278	54309	0.378	ng/ul#	96
26) Benzo(g,h,i)perylene	25.94	276	55424	0.367	ng/ul	97

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

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