

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007470.D
 Acq On : 28 Aug 2019 16:13
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.436

Quant Time: Aug 28 16:59:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 20:30:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	4039	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	17730	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	9093	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	20205	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	19039	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	19729	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	10781	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	26839	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.20	128	19423	0.388	ng/ul	99
5) 2-Methylnaphthalene	11.83	142	13182	0.392	ng/ul	99
7) Acenaphthylene	13.76	152	19311	0.382	ng/ul	99
8) Acenaphthene	14.11	153	14122	0.388	ng/ul	98
9) Fluorene	15.11	166	15254	0.388	ng/ul	98
11) Pentachlorophenol	16.47	266	1916	0.421	ng/ul	99
12) Phenanthrene	16.84	178	25240	0.400	ng/ul	99
13) Anthracene	16.94	178	24257	0.408	ng/ul	100
15) Fluoranthene	18.87	202	33993	0.403	ng/ul	99
17) Pyrene	19.24	202	35201	0.378	ng/ul	99
18) Benzo(a)anthracene	21.01	228	33152	0.389	ng/ul	100
19) Chrysene	21.06	228	34906	0.382	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	33602	0.390	ng/ul	99
22) Benzo(k)fluoranthene	22.55	252	35686	0.396	ng/ul#	93
23) Benzo(a)pyrene	23.03	252	33106	0.385	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.20	276	38409	0.400	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	30069	0.395	ng/ul	100
26) Benzo(g,h,i)perylene	25.81	276	32016	0.390	ng/ul	99

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

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