

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007491.D
 Acq On : 29 Aug 2019 05:21
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.438

Quant Time: Aug 29 06:42:36 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 : Thu Aug 29 04:22:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	4448	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	19129	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	9808	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	21342	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	18450	0.40	ng/ul	0.00
20) Perylene-d12	23.11	264	19120	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	11726	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	26409	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	21186	0.392	ng/ul	99
5) 2-Methylnaphthalene	11.83	142	14277	0.393	ng/ul	99
7) Acenaphthylene	13.76	152	20386	0.374	ng/ul	99
8) Acenaphthene	14.11	153	15230	0.388	ng/ul	99
9) Fluorene	15.11	166	16271	0.383	ng/ul	97
11) Pentachlorophenol	16.47	266	1819	0.378	ng/ul	99
12) Phenanthrene	16.84	178	26692	0.401	ng/ul	99
13) Anthracene	16.93	178	25076	0.400	ng/ul	99
15) Fluoranthene	18.87	202	33388	0.375	ng/ul	100
17) Pyrene	19.24	202	34295	0.380	ng/ul	99
18) Benzo(a)anthracene	21.00	228	31311	0.380	ng/ul	99
19) Chrysene	21.05	228	33615	0.380	ng/ul	99
21) Benzo(b)fluoranthene	22.50	252	32023	0.383	ng/ul	99
22) Benzo(k)fluoranthene	22.54	252	34805	0.399	ng/ul#	95
23) Benzo(a)pyrene	23.03	252	31607	0.380	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.18	276	36760	0.395	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.19	278	29045	0.394	ng/ul	100
26) Benzo(g,h,i)perylene	25.81	276	30827	0.387	ng/ul	99

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

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