

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072219\
 Data File : BN006960.D
 Acq On : 23 Jul 2019 02:21
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.443

Quant Time: Jul 23 04:28:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 23 04:26:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1890	0.40	ng/ul	0.01
2) Naphthalene-d8	10.38	136	7659	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4035	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8226	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6179	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	6733	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	4594	0.39	ng/ul	0.01
14) Fluoranthene-d10	19.06	212	9024	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.43	128	9218	0.419	ng/ul	97
5) 2-Methylnaphthalene	12.07	142	5465	0.368	ng/ul	96
7) Acenaphthylene	13.97	152	7773	0.373	ng/ul	99
8) Acenaphthene	14.31	153	6540	0.407	ng/ul	98
9) Fluorene	15.32	166	6569	0.373	ng/ul	99
11) Pentachlorophenol	16.73	266	357	0.291	ng/ul#	91
12) Phenanthrene	17.06	178	10424	0.394	ng/ul	98
13) Anthracene	17.17	178	7651	0.335	ng/ul	98
15) Fluoranthene	19.09	202	11811	0.361	ng/ul	100
17) Pyrene	19.46	202	12549	0.444	ng/ul	99
18) Benzo(a)anthracene	21.22	228	8454	0.333	ng/ul	99
19) Chrysene	21.27	228	13422	0.474	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	11029	0.400	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	12756	0.442	ng/ul	98
23) Benzo(a)pyrene	23.37	252	11198	0.414	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.70	276	12792	0.468	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.70	278	9803	0.451	ng/ul	97
26) Benzo(g,h,i)perylene	26.38	276	11634	0.509	ng/ul	98

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

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