

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072219\
 Data File : BN006945.D
 Acq On : 22 Jul 2019 17:40
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.442

Quant Time: Jul 23 04:25:12 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 04:22:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	2211	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	8275	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4207	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8636	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	6153	0.40	ng/ul	0.01
20) Perylene-d12	23.49	264	6855	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	5062	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	9446	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	10066	0.424	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	6179	0.385	ng/ul	99
7) Acenaphthylene	13.97	152	8149	0.375	ng/ul	99
8) Acenaphthene	14.31	153	6881	0.411	ng/ul	98
9) Fluorene	15.32	166	7049	0.384	ng/ul	100
11) Pentachlorophenol	16.73	266	550	0.428	ng/ul	97
12) Phenanthrene	17.06	178	11171	0.402	ng/ul	99
13) Anthracene	17.17	178	8085	0.337	ng/ul	98
15) Fluoranthene	19.09	202	12502	0.364	ng/ul	99
17) Pyrene	19.46	202	13245	0.471	ng/ul	100
18) Benzo(a)anthracene	21.24	228	8692	0.344	ng/ul	100
19) Chrysene	21.29	228	13838	0.491	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	11105	0.395	ng/ul	98
22) Benzo(k)fluoranthene	22.86	252	12849	0.437	ng/ul	99
23) Benzo(a)pyrene	23.39	252	10914	0.397	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.73	276	12690	0.456	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.73	278	8947	0.404	ng/ul	97
26) Benzo(g,h,i)perylene	26.41	276	11712	0.503	ng/ul	98

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

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