

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071119\
 Data File : BN006837.D
 Acq On : 12 Jul 2019 16:54
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.426

Quant Time: Jul 12 17:27:41 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 : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2630	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	9930	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	5221	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	11170	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	9008	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	10220	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	5718	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	11862	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	11149	0.391	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	6982	0.362	ng/ul	99
7) Acenaphthylene	13.96	152	9851	0.366	ng/ul	99
8) Acenaphthene	14.30	153	7855	0.378	ng/ul	99
9) Fluorene	15.31	166	8271	0.363	ng/ul	100
11) Pentachlorophenol	16.71	266	425	0.255	ng/ul#	85
12) Phenanthrene	17.05	178	13260	0.369	ng/ul	100
13) Anthracene	17.15	178	10858	0.350	ng/ul	99
15) Fluoranthene	19.08	202	15292	0.344	ng/ul	99
17) Pyrene	19.45	202	16104	0.391	ng/ul	98
18) Benzo(a)anthracene	21.22	228	13038	0.352	ng/ul	100
19) Chrysene	21.27	228	16099	0.390	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	15052	0.359	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	16862	0.385	ng/ul	99
23) Benzo(a)pyrene	23.37	252	15070	0.367	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.69	276	17948	0.432	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.69	278	14101	0.427	ng/ul	98
26) Benzo(g,h,i)perylene	26.37	276	15645	0.451	ng/ul	96

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

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