

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011381.D
 Acq On : 06 Aug 2020 16:14
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Aug 06 17:01:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	878	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	2772	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1662	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3588	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3431	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3573	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	2570	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	6122	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	3322	0.403	ng/ul	95
5) 2-Methylnaphthalene	11.88	142	2019	0.365	ng/ul	99
7) Acenaphthylene	13.81	152	3012	0.379	ng/ul	98
8) Acenaphthene	14.15	153	2275	0.361	ng/ul	98
9) Fluorene	15.15	166	2655	0.358	ng/ul	99
11) Pentachlorophenol	16.52	266	310	0.352	ng/ul	98
12) Phenanthrene	16.89	178	4167	0.351	ng/ul	96
13) Anthracene	16.98	178	3362	0.324	ng/ul	95
15) Fluoranthene	18.92	202	5115	0.349	ng/ul	97
17) Pyrene	19.28	202	4679	0.300	ng/ul	99
18) Benzo(a)anthracene	21.05	228	4398	0.334	ng/ul	98
19) Chrysene	21.10	228	4928	0.326	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	4560	0.306	ng/ul	92
22) Benzo(k)fluoranthene	22.60	252	4794	0.287	ng/ul	94
23) Benzo(a)pyrene	23.09	252	4829	0.372	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.25	276	5831	0.428	ng/ul	97
25) Dibenzo(a,h)anthracene	25.26	278	4696	0.433	ng/ul	96
26) Benzo(g,h,i)perylene	25.86	276	4640	0.377	ng/ul	97

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

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