

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080320\
 Data File : BN011346.D
 Acq On : 03 Aug 2020 23:03
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Aug 04 06:12:24 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 : Mon Aug 03 10:06:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	1211	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4744	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	3133	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	6590	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	6048	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	6796	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	5404	0.45	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	13915	0.44	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.26	128	5879	0.417	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	4480	0.473	ng/ul	99
7) Acenaphthylene	13.81	152	6672	0.446	ng/ul	98
8) Acenaphthene	14.15	153	4727	0.398	ng/ul	100
9) Fluorene	15.15	166	5463	0.391	ng/ul	99
11) Pentachlorophenol	16.51	266	728	0.451	ng/ul	97
12) Phenanthrene	16.89	178	9076	0.416	ng/ul	99
13) Anthracene	16.98	178	8037	0.422	ng/ul	98
15) Fluoranthene	18.92	202	11701	0.435	ng/ul	98
17) Pyrene	19.28	202	11539	0.420	ng/ul	100
18) Benzo(a)anthracene	21.05	228	9827	0.423	ng/ul	100
19) Chrysene	21.10	228	11264	0.423	ng/ul	100
21) Benzo(b)fluoranthene	22.56	252	11445	0.403	ng/ul	100
22) Benzo(k)fluoranthene	22.60	252	13545	0.426	ng/ul	97
23) Benzo(a)pyrene	23.09	252	10374	0.421	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.25	276	10851	0.419	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.26	278	8684	0.421	ng/ul	99
26) Benzo(g,h,i)perylene	25.87	276	9326	0.398	ng/ul	100

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

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