

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080320\
 Data File : BN011334.D
 Acq On : 03 Aug 2020 15:51
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.414

Quant Time: Aug 03 16:22:09 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	7.47	152	1085	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	3912	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2397	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	5427	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	5059	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	5198	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	4127	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	10640	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.26	128	4817	0.414	ng/ul	98
5) 2-Methylnaphthalene	11.89	142	3261	0.418	ng/ul	98
7) Acenaphthylene	13.81	152	4834	0.422	ng/ul	99
8) Acenaphthene	14.16	153	3878	0.426	ng/ul	99
9) Fluorene	15.16	166	4478	0.419	ng/ul	99
11) Pentachlorophenol	16.52	266	529	0.398	ng/ul	98
12) Phenanthrene	16.89	178	7533	0.419	ng/ul	99
13) Anthracene	16.99	178	6558	0.418	ng/ul	99
15) Fluoranthene	18.92	202	9054	0.408	ng/ul	98
17) Pyrene	19.29	202	9069	0.395	ng/ul	99
18) Benzo(a)anthracene	21.05	228	7938	0.408	ng/ul	100
19) Chrysene	21.10	228	9570	0.429	ng/ul	100
21) Benzo(b)fluoranthene	22.57	252	8760	0.404	ng/ul	99
22) Benzo(k)fluoranthene	22.61	252	10265	0.422	ng/ul	99
23) Benzo(a)pyrene	23.10	252	7900	0.419	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.26	276	8937	0.451	ng/ul	97
25) Dibenzo(a,h)anthracene	25.28	278	7098	0.450	ng/ul	98
26) Benzo(g,h,i)perylene	25.88	276	7694	0.430	ng/ul	99

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

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