

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011712.D
 Acq On : 04 Sep 2020 22:22
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Sep 05 05:45:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 05 01:53:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1821	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	6764	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3746	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	8171	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	6772	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	6322	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	4404	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	8826	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	7931	0.402	ng/ul	98
5) 2-Methylnaphthalene	12.19	142	5370	0.431	ng/ul	98
7) Acenaphthylene	14.10	152	7193	0.381	ng/ul	97
8) Acenaphthene	14.44	153	5809	0.396	ng/ul	99
9) Fluorene	15.43	166	6664	0.385	ng/ul	99
11) Pentachlorophenol	16.77	266	912	0.470	ng/ul	99
12) Phenanthrene	17.16	178	10979	0.397	ng/ul	97
13) Anthracene	17.26	178	9811	0.431	ng/ul	97
15) Fluoranthene	19.19	202	12213	0.384	ng/ul	99
17) Pyrene	19.55	202	12271	0.464	ng/ul	98
18) Benzo(a)anthracene	21.30	228	10608	0.467	ng/ul	98
19) Chrysene	21.35	228	11878	0.388	ng/ul	100
21) Benzo(b)fluoranthene	22.89	252	11724	0.427	ng/ul	96
22) Benzo(k)fluoranthene	22.94	252	12439	0.409	ng/ul	97
23) Benzo(a)pyrene	23.47	252	10339	0.448	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.83	276	14059	0.450	ng/ul	96
25) Dibenzo(a,h)anthracene	25.85	278	11479	0.451	ng/ul	97
26) Benzo(g,h,i)perylene	26.52	276	11814	0.404	ng/ul	98

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

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