

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011739.D
 Acq On : 08 Sep 2020 16:30
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.497

Quant Time: Sep 08 17:02:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 14:34:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1666	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	5951	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3204	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	6944	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5678	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	5236	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	4038	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7771	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	7642	0.434	ng/ul	100
5) 2-Methylnaphthalene	12.19	142	5112	0.435	ng/ul	100
7) Acenaphthylene	14.09	152	6737	0.436	ng/ul	99
8) Acenaphthene	14.44	153	5408	0.436	ng/ul	99
9) Fluorene	15.43	166	6248	0.436	ng/ul	99
11) Pentachlorophenol	16.77	266	880	0.420	ng/ul	100
12) Phenanthrene	17.16	178	10194	0.436	ng/ul	98
13) Anthracene	17.25	178	9028	0.434	ng/ul	99
15) Fluoranthene	19.18	202	11208	0.432	ng/ul	99
17) Pyrene	19.54	202	11129	0.441	ng/ul	98
18) Benzo(a)anthracene	21.30	228	9560	0.433	ng/ul	99
19) Chrysene	21.35	228	10708	0.434	ng/ul	98
21) Benzo(b)fluoranthene	22.89	252	10664	0.435	ng/ul	98
22) Benzo(k)fluoranthene	22.93	252	11058	0.439	ng/ul	100
23) Benzo(a)pyrene	23.46	252	9340	0.441	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.82	276	12166	0.437	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	9827	0.432	ng/ul	100
26) Benzo(g,h,i)perylene	26.51	276	10247	0.421	ng/ul	98

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

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