

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008463.D
 Acq On : 06 Nov 2019 00:47
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.424

Quant Time: Nov 07 11:11:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 10:57:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1323	0.40	ng/ul	0.01
2) Naphthalene-d8	10.32	136	5895	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3712	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6947	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	4958	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	4453	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.93	152	3581	0.38	ng/ul	0.01
14) Fluoranthene-d10	18.99	212	8290	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.37	128	6678	0.398	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	4306	0.385	ng/ul	98
7) Acenaphthylene	13.92	152	6262	0.362	ng/ul	100
8) Acenaphthene	14.26	153	5228	0.373	ng/ul	99
9) Fluorene	15.26	166	5551	0.362	ng/ul	100
11) Pentachlorophenol	16.63	266	516	0.341	ng/ul	98
12) Phenanthrene	16.99	178	8378	0.397	ng/ul	100
13) Anthracene	17.09	178	7114	0.385	ng/ul	100
15) Fluoranthene	19.02	202	9675	0.400	ng/ul	99
17) Pyrene	19.38	202	10117	0.418	ng/ul	98
18) Benzo(a)anthracene	21.14	228	6291	0.364	ng/ul	99
19) Chrysene	21.20	228	9219	0.420	ng/ul	99
21) Benzo(b)fluoranthene	22.68	252	6665	0.402	ng/ul	93
22) Benzo(k)fluoranthene	22.72	252	7643	0.405	ng/ul	99
23) Benzo(a)pyrene	23.24	252	6696	0.401	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.47	276	7552	0.398	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.49	278	5940	0.395	ng/ul	99
26) Benzo(g,h,i)perylene	26.12	276	6971	0.410	ng/ul	99

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

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