

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008441.D
 Acq On : 05 Nov 2019 11:17
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Nov 05 14:03:30 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 : Tue Nov 05 13:48:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1381	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5918	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3662	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	7037	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	4969	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	4584	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	3714	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	8275	0.34	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.37	128	6757	0.408	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	4457	0.387	ng/ul	99
7) Acenaphthylene	13.92	152	6299	0.365	ng/ul	100
8) Acenaphthene	14.26	153	5140	0.383	ng/ul	99
9) Fluorene	15.25	166	5622	0.367	ng/ul	100
11) Pentachlorophenol	16.62	266	610	0.390	ng/ul	97
12) Phenanthrene	16.99	178	8553	0.417	ng/ul	99
13) Anthracene	17.08	178	7474	0.408	ng/ul	99
15) Fluoranthene	19.02	202	9651	0.358	ng/ul	99
17) Pyrene	19.39	202	10008	0.518	ng/ul	99
18) Benzo(a)anthracene	21.15	228	7043	0.392	ng/ul	99
19) Chrysene	21.20	228	8795	0.404	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	6388	0.427	ng/ul	93
22) Benzo(k)fluoranthene	22.73	252	7560	0.451	ng/ul	97
23) Benzo(a)pyrene	23.24	252	6673	0.429	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.48	276	7775	0.421	ng/ul	98
25) Dibenzo(a,h)anthracene	25.49	278	6232	0.425	ng/ul	96
26) Benzo(g,h,i)perylene	26.13	276	6892	0.410	ng/ul	99

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

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