

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN051719\
 Data File : BN005742.D
 Acq On : 17 May 2019 22:27
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.481

Quant Time: May 17 23:18:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 17 04:47:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1977	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7884	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	4694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	10235	0.40	ng/ul	0.01
16) Chrysene-d12	21.23	240	7433	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7648	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5289	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	11206	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	8527	0.409	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	6151	0.432	ng/ul	99
7) Acenaphthylene	13.95	152	10246	0.434	ng/ul	99
8) Acenaphthene	14.30	153	6930	0.414	ng/ul	98
9) Fluorene	15.30	166	8051	0.412	ng/ul	99
11) Pentachlorophenol	16.70	266	851	0.321	ng/ul	91
12) Phenanthrene	17.04	178	11862	0.425	ng/ul	99
13) Anthracene	17.13	178	11776	0.423	ng/ul	99
15) Fluoranthene	19.07	202	13283	0.344	ng/ul	96
17) Pyrene	19.44	202	13410	0.489	ng/ul#	95
18) Benzo(a)anthracene	21.21	228	11402	0.445	ng/ul#	92
19) Chrysene	21.27	228	11691	0.371	ng/ul#	90
21) Benzo(b)fluoranthene	22.79	252	10858	0.472	ng/ul	81
22) Benzo(k)fluoranthene	22.83	252	11308	0.392	ng/ul#	82
23) Benzo(a)pyrene	23.35	252	10830	0.421	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.66	276	12611	0.439	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.66	278	10202	0.429	ng/ul#	82
26) Benzo(g,h,i)perylene	26.34	276	10763	0.439	ng/ul#	88

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

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