

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005492.D
 Acq On : 06 May 2019 19:19
 Operator : JU/SJ
 Sample : SSTD0.861
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.861

Quant Time: May 07 02:51:19 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 : Tue May 07 02:47:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1929	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7350	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	4248	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	10501	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	11396	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	12941	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	9496	0.81	ng/ul	-0.01
14) Fluoranthene-d10	19.03	212	26541	0.88	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.41	128	15528	0.775	ng/ul	97
5) 2-Methylnaphthalene	12.04	142	11029	0.780	ng/ul	100
7) Acenaphthylene	13.95	152	17412	0.816	ng/ul	98
8) Acenaphthene	14.30	153	12199	0.776	ng/ul	97
9) Fluorene	15.30	166	14327	0.788	ng/ul	98
11) Pentachlorophenol	16.66	266	2158	0.847	ng/ul	93
12) Phenanthrene	17.04	178	23288	0.731	ng/ul	100
13) Anthracene	17.13	178	23415	0.795	ng/ul	98
15) Fluoranthene	19.06	202	32011	0.857	ng/ul	99
17) Pyrene	19.43	202	33271	0.700	ng/ul	98
18) Benzo(a)anthracene	21.21	228	32622	0.765	ng/ul	99
19) Chrysene	21.26	228	36974	0.815	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	33189	0.698	ng/ul	97
22) Benzo(k)fluoranthene	22.82	252	38834	0.791	ng/ul	97
23) Benzo(a)pyrene	23.34	252	35169	0.801	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.63	276	39137	0.803	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.64	278	32250	0.807	ng/ul	97
26) Benzo(g,h,i)perylene	26.28	276	32576	0.786	ng/ul	97

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

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