

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053019\
 Data File : BN005957.D
 Acq On : 30 May 2019 13:50
 Operator : JU/SJ
 Sample : SSTD0.891
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.891

Quant Time: May 30 15:08:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 15:01:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	8138	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	30139	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	16487	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.15	188	38916	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	29395	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	26236	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	37730	0.83	ng/ul	0.00
14) Fluoranthene-d10	19.18	212	88122	0.87	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.58	128	67259	0.815	ng/ul	99
5) 2-Methylnaphthalene	12.20	142	47367	0.830	ng/ul	100
7) Acenaphthylene	14.11	152	74433	0.825	ng/ul	100
8) Acenaphthene	14.45	153	51562	0.820	ng/ul	100
9) Fluorene	15.44	166	60635	0.826	ng/ul	98
11) Pentachlorophenol	16.79	266	7973	1.016	ng/ul	99
12) Phenanthrene	17.18	178	100747	0.813	ng/ul	100
13) Anthracene	17.28	178	97718	0.839	ng/ul	100
15) Fluoranthene	19.21	202	115372	0.856	ng/ul	98
17) Pyrene	19.57	202	116936	0.744	ng/ul	99
18) Benzo(a)anthracene	21.34	228	104349	0.807	ng/ul	98
19) Chrysene	21.39	228	97777	0.803	ng/ul	99
21) Benzo(b)fluoranthene	22.95	252	89763	0.833	ng/ul	93
22) Benzo(k)fluoranthene	22.99	252	87025	0.826	ng/ul	94
23) Benzo(a)pyrene	23.53	252	82554	0.817	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.92	276	84394	0.753	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.94	278	67152	0.748	ng/ul	95
26) Benzo(g,h,i)perylene	26.62	276	69566	0.738	ng/ul	97

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

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