

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005145.D
 Acq On : 18 Apr 2019 09:03
 Operator : JU/SJ
 Sample : SSTD0.253
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.253

Quant Time: Apr 18 11:18:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 11:12:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	3953	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	14353	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7878	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	18485	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	14849	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	14439	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4448	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	9938	0.17	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	7761	0.188	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	5426	0.184	ng/ul	100
7) Acenaphthylene	13.95	152	7703	0.214	ng/ul	100
8) Acenaphthene	14.30	153	5810	0.186	ng/ul	99
9) Fluorene	15.30	166	6667	0.178	ng/ul	99
11) Pentachlorophenol	16.64	266	713	0.181	ng/ul	98
12) Phenanthrene	17.03	178	11185	0.186	ng/ul	99
13) Anthracene	17.12	178	10090	0.199	ng/ul	99
15) Fluoranthene	19.06	202	12381	0.168	ng/ul	99
17) Pyrene	19.42	202	12436	0.260	ng/ul	97
18) Benzo(a)anthracene	21.19	228	10907	0.228	ng/ul	99
19) Chrysene	21.24	228	11517	0.207	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	10413	0.180	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	10338	0.182	ng/ul	98
23) Benzo(a)pyrene	23.29	252	9288	0.180	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.56	276	10438	0.154	ng/ul	99
25) Dibenzo(a,h)anthracene	25.58	278	8570	0.156	ng/ul	96
26) Benzo(g,h,i)perylene	26.22	276	9011	0.146	ng/ul	98

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

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