

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005146.D
 Acq On : 18 Apr 2019 09:38
 Operator : JU/SJ
 Sample : SSTD0.454
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.454

Quant Time: Apr 18 11:17:07 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	3044	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	11316	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	6429	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	14818	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	11826	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	11108	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	7421	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	16630	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.42	128	12556	0.386	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	8941	0.385	ng/ul	100
7) Acenaphthylene	13.95	152	12733	0.433	ng/ul	100
8) Acenaphthene	14.30	153	9681	0.380	ng/ul	100
9) Fluorene	15.30	166	11230	0.368	ng/ul	100
11) Pentachlorophenol	16.64	266	1286	0.408	ng/ul	100
12) Phenanthrene	17.03	178	18581	0.386	ng/ul	100
13) Anthracene	17.12	178	16804	0.414	ng/ul	100
15) Fluoranthene	19.06	202	20804	0.352	ng/ul	100
17) Pyrene	19.42	202	20593	0.541	ng/ul	100
18) Benzo(a)anthracene	21.19	228	17948	0.471	ng/ul	100
19) Chrysene	21.24	228	18662	0.422	ng/ul	100
21) Benzo(b)fluoranthene	22.74	252	16858	0.380	ng/ul	100
22) Benzo(k)fluoranthene	22.78	252	16704	0.383	ng/ul	100
23) Benzo(a)pyrene	23.29	252	15071	0.379	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.56	276	16636	0.319	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	13595	0.321	ng/ul	100
26) Benzo(g,h,i)perylene	26.21	276	14318	0.302	ng/ul	100

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

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