

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002500.D
 Acq On : 24 Aug 2018 13:34
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Aug 24 15:06:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 24 02:20:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1156	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	5253	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3281	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	8553	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9325	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	9802	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	3474	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11321	0.45	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	5455	0.403	ng/ul#	96
5) 2-Methylnaphthalene	12.13	142	3941	0.418	ng/ul	98
7) Acenaphthylene	14.03	152	6513	0.424	ng/ul	97
8) Acenaphthene	14.37	153	4770	0.407	ng/ul	95
9) Fluorene	15.37	166	5547	0.413	ng/ul	92
11) Pentachlorophenol	16.74	266	548	0.475	ng/ul	98
12) Phenanthrene	17.10	178	9960	0.394	ng/ul#	90
13) Anthracene	17.20	178	9907	0.456	ng/ul#	93
15) Fluoranthene	19.12	202	12720	0.428	ng/ul	97
17) Pyrene	19.49	202	13180	0.414	ng/ul	96
18) Benzo(a)anthracene	21.24	228	12782	0.458	ng/ul#	87
19) Chrysene	21.29	228	12608	0.400	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	13213	0.375	ng/ul#	56
22) Benzo(k)fluoranthene	22.86	252	13924	0.383	ng/ul#	57
23) Benzo(a)pyrene	23.38	252	13113	0.396	ng/ul#	42
24) Indeno(1,2,3-cd)pyrene	25.70	276	14750	0.428	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.70	278	11749	0.432	ng/ul#	73
26) Benzo(g,h,i)perylene	26.37	276	12528	0.424	ng/ul#	94

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

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