

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092218\
 Data File : BN002809.D
 Acq On : 21 Sep 2018 08:52
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.471

Quant Time: Sep 21 22:59:56 2018
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SOM-EPA-SIM-BN092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 21 04:00:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1366	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	5988	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	3284	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	6966	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	5484	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	5193	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	3641	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	7067	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.48	128	6172	0.410	ng/ul#	97
5) 2-Methylnaphthalene	12.12	142	4163	0.426	ng/ul	99
7) Acenaphthylene	14.01	152	5606	0.404	ng/ul	96
8) Acenaphthene	14.36	153	4679	0.413	ng/ul	97
9) Fluorene	15.36	166	5124	0.424	ng/ul	93
11) Pentachlorophenol	16.75	266	222	0.283	ng/ul	98
12) Phenanthrene	17.09	178	7917	0.412	ng/ul#	90
13) Anthracene	17.20	178	6811	0.408	ng/ul#	92
15) Fluoranthene	19.12	202	8144	0.393	ng/ul	96
17) Pyrene	19.49	202	8329	0.466	ng/ul	98
18) Benzo(a)anthracene	21.24	228	6475	0.427	ng/ul#	85
19) Chrysene	21.29	228	7661	0.433	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	6961	0.406	ng/ul	89
22) Benzo(k)fluoranthene	22.85	252	7606	0.421	ng/ul#	91
23) Benzo(a)pyrene	23.38	252	6746	0.410	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.69	276	7646	0.386	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.70	278	6083	0.383	ng/ul	94
26) Benzo(g,h,i)perylene	26.37	276	6685	0.391	ng/ul	98

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

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