

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092218\
 Data File : BN002821.D
 Acq On : 21 Sep 2018 16:58
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.472

Quant Time: Sep 21 23:10:55 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 23:09:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	1226	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	5219	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	2718	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	5548	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4994	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	4830	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	3121	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6129	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.48	128	5406	0.412	ng/ul#	97
5) 2-Methylnaphthalene	12.12	142	3592	0.422	ng/ul	99
7) Acenaphthylene	14.01	152	4689	0.409	ng/ul	99
8) Acenaphthene	14.36	153	3847	0.411	ng/ul	97
9) Fluorene	15.36	166	4220	0.422	ng/ul	93
11) Pentachlorophenol	16.74	266	290	0.464	ng/ul	97
12) Phenanthrene	17.09	178	6434	0.420	ng/ul#	89
13) Anthracene	17.20	178	5716	0.430	ng/ul#	93
15) Fluoranthene	19.12	202	6985	0.423	ng/ul	95
17) Pyrene	19.49	202	7108	0.436	ng/ul	98
18) Benzo(a)anthracene	21.24	228	5977	0.433	ng/ul#	86
19) Chrysene	21.29	228	6839	0.425	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	6670	0.418	ng/ul	89
22) Benzo(k)fluoranthene	22.86	252	7062	0.421	ng/ul#	92
23) Benzo(a)pyrene	23.39	252	6386	0.418	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.70	276	7518	0.408	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.71	278	6008	0.407	ng/ul	96
26) Benzo(g,h,i)perylene	26.38	276	6542	0.411	ng/ul	98

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

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