

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097289.D
 Acq On : 17 Aug 2018 9:45
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.401

Quant Time: Aug 17 12:19:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 12:19:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	966	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5103	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2743	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	6241	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	5280	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4436	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3077	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5996	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	4635	0.375	ng/ul#	92
5) 2-Methylnaphthalene	11.80	142	3014	0.369	ng/ul	100
7) Acenaphthylene	13.73	152	4584	0.386	ng/ul	98
8) Acenaphthene	14.08	153	3303	0.385	ng/ul#	89
9) Fluorene	15.08	166	3569	0.362	ng/ul#	92
11) Pentachlorophenol	16.45	266	132	0.148	ng/ul	95
12) Phenanthrene	16.81	178	5928	0.373	ng/ul#	88
13) Anthracene	16.90	178	4981	0.336	ng/ul#	91
15) Fluoranthene	18.84	202	6257	0.298	ng/ul	85
17) Pyrene	19.21	202	5418	0.383	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	4679	0.310	ng/ul#	84
19) Chrysene	21.02	228	5866	0.377	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	4324	0.358	ng/ul	86
22) Benzo(k)fluoranthene	22.59	252	5530	0.424	ng/ul#	87
23) Benzo(a)pyrene	23.12	252	4147	0.343	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.53	276	4886	0.323	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.55	278	3930	0.317	ng/ul	95
26) Benzo(g,h,i)perylene	26.22	276	4119	0.315	ng/ul	93

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

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