

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097300.D
 Acq On : 17 Aug 2018 17:10
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.408

Quant Time: Aug 17 23:15:31 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 14:15:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	948	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4956	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2578	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	5790	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	4861	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4020	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2964	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5671	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	4518	0.374	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	2901	0.372	ng/ul	99
7) Acenaphthylene	13.73	152	4284	0.368	ng/ul	98
8) Acenaphthene	14.08	153	3134	0.374	ng/ul#	88
9) Fluorene	15.08	166	3364	0.369	ng/ul	92
11) Pentachlorophenol	16.45	266	103	0.274	ng/ul	90
12) Phenanthrene	16.81	178	5518	0.369	ng/ul#	89
13) Anthracene	16.90	178	4577	0.365	ng/ul#	92
15) Fluoranthene	18.84	202	5815	0.366	ng/ul	86
17) Pyrene	19.21	202	5023	0.342	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	4382	0.364	ng/ul#	85
19) Chrysene	21.02	228	5442	0.378	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	4082	0.375	ng/ul	86
22) Benzo(k)fluoranthene	22.58	252	5071	0.375	ng/ul#	89
23) Benzo(a)pyrene	23.12	252	3793	0.377	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.53	276	4457	0.356	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.55	278	3621	0.358	ng/ul	94
26) Benzo(g,h,i)perylene	26.22	276	3698	0.369	ng/ul	93

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

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