

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE082418\
 Data File : BE097359.D
 Acq On : 24 Aug 2018 8:58
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.411

Quant Time: Aug 24 22:48:19 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 : Sat Aug 18 05:45:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	800	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3768	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2025	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	5681	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	7297	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	6709	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2359	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	7603	0.50	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.17	128	3719	0.405	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	2353	0.397	ng/ul	98
7) Acenaphthylene	13.73	152	3015	0.330	ng/ul	99
8) Acenaphthene	14.07	153	2592	0.394	ng/ul	95
9) Fluorene	15.07	166	2984	0.417	ng/ul#	92
11) Pentachlorophenol	16.44	266	201	0.545	ng/ul	98
12) Phenanthrene	16.80	178	5929	0.404	ng/ul#	88
13) Anthracene	16.90	178	4946	0.402	ng/ul#	92
15) Fluoranthene	18.84	202	7600	0.487	ng/ul	84
17) Pyrene	19.20	202	7383	0.335	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	7615	0.421	ng/ul#	86
19) Chrysene	21.01	228	8868	0.411	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	8110	0.446	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	7884	0.350	ng/ul#	86
23) Benzo(a)pyrene	23.10	252	7016	0.418	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.50	276	8981	0.430	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.52	278	7436	0.440	ng/ul#	89
26) Benzo(g,h,i)perylene	26.19	276	7905	0.473	ng/ul#	91

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

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