

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097133.D
 Acq On : 1 Aug 2018 17:29
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.480

Quant Time: Aug 02 01:03:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 01 16:16:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1424	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6633	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3632	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9938	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12106	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11745	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4260	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12870	0.38	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.18	128	6436	0.401	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	4221	0.398	ng/ul	98
7) Acenaphthylene	13.74	152	6245	0.398	ng/ul	97
8) Acenaphthene	14.08	153	4563	0.402	ng/ul#	94
9) Fluorene	15.08	166	5177	0.396	ng/ul#	94
11) Pentachlorophenol	16.44	266	461	0.324	ng/ul	96
12) Phenanthrene	16.81	178	9988	0.394	ng/ul#	88
13) Anthracene	16.90	178	9126	0.386	ng/ul#	91
15) Fluoranthene	18.84	202	12727	0.380	ng/ul	84
17) Pyrene	19.21	202	13149	0.405	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	13291	0.384	ng/ul#	84
19) Chrysene	21.01	228	14083	0.395	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	12343	0.386	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	13710	0.397	ng/ul#	87
23) Benzo(a)pyrene	23.11	252	11943	0.373	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	15256	0.381	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.53	278	12631	0.385	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	13461	0.389	ng/ul#	92

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

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