

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097226.D
 Acq On : 8 Aug 2018 9:24
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.493

Quant Time: Aug 08 23:24:44 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 : Mon Aug 06 15:47:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1217	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6666	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3843	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	8849	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	8718	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8808	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4301	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	10077	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	6209	0.385	ng/ul#	92
5) 2-Methylnaphthalene	11.80	142	4262	0.400	ng/ul	98
7) Acenaphthylene	13.73	152	6601	0.397	ng/ul	96
8) Acenaphthene	14.08	153	4657	0.388	ng/ul#	94
9) Fluorene	15.08	166	5147	0.373	ng/ul#	93
11) Pentachlorophenol	16.45	266	366	0.289	ng/ul	97
12) Phenanthrene	16.81	178	8804	0.390	ng/ul#	89
13) Anthracene	16.90	178	8220	0.391	ng/ul#	92
15) Fluoranthene	18.84	202	9958	0.334	ng/ul	84
17) Pyrene	19.21	202	9760	0.418	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	9543	0.383	ng/ul#	84
19) Chrysene	21.01	228	10029	0.391	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	9241	0.385	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	10201	0.394	ng/ul#	86
23) Benzo(a)pyrene	23.11	252	9475	0.394	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	11874	0.395	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.53	278	9742	0.396	ng/ul#	90
26) Benzo(g,h,i)perylene	26.21	276	10040	0.386	ng/ul#	93

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

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