

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097241.D
 Acq On : 8 Aug 2018 18:51
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.494

Quant Time: Aug 08 23:27:37 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 08 23:25:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1177	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6297	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3566	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	8190	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	8300	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8576	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	4057	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9210	0.33	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.18	128	5905	0.387	ng/ul#	92
5) 2-Methylnaphthalene	11.80	142	4029	0.400	ng/ul	97
7) Acenaphthylene	13.73	152	6175	0.400	ng/ul	97
8) Acenaphthene	14.08	153	4411	0.396	ng/ul	94
9) Fluorene	15.08	166	4888	0.381	ng/ul#	95
11) Pentachlorophenol	16.44	266	386	0.329	ng/ul	96
12) Phenanthrene	16.81	178	8024	0.384	ng/ul#	89
13) Anthracene	16.90	178	7675	0.394	ng/ul#	92
15) Fluoranthene	18.84	202	9152	0.332	ng/ul	84
17) Pyrene	19.21	202	9224	0.415	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	9454	0.399	ng/ul#	85
19) Chrysene	21.01	228	9451	0.387	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	9689	0.415	ng/ul	85
22) Benzo(k)fluoranthene	22.58	252	9312	0.369	ng/ul#	89
23) Benzo(a)pyrene	23.11	252	9251	0.396	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.51	276	11704	0.400	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.53	278	9792	0.409	ng/ul#	90
26) Benzo(g,h,i)perylene	26.21	276	9843	0.389	ng/ul#	93

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

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