

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097154.D
 Acq On : 2 Aug 2018 16:55
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.483

Quant Time: Aug 02 18:13:10 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 : Thu Aug 02 14:49:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1261	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5964	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3336	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9138	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10484	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10306	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3805	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	11666	0.37	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	5850	0.405	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	3787	0.397	ng/ul	98
7) Acenaphthylene	13.73	152	5808	0.403	ng/ul	97
8) Acenaphthene	14.08	153	4170	0.400	ng/ul#	93
9) Fluorene	15.08	166	4766	0.397	ng/ul#	94
11) Pentachlorophenol	16.44	266	454	0.347	ng/ul	95
12) Phenanthrene	16.81	178	9262	0.398	ng/ul#	88
13) Anthracene	16.90	178	8555	0.394	ng/ul#	91
15) Fluoranthene	18.84	202	11666	0.379	ng/ul	84
17) Pyrene	19.20	202	12163	0.433	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	11399	0.380	ng/ul#	85
19) Chrysene	21.01	228	12213	0.396	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	10570	0.376	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	11963	0.395	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	10297	0.366	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	13129	0.373	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.53	278	11092	0.385	ng/ul	91
26) Benzo(g,h,i)perylene	26.20	276	11302	0.372	ng/ul#	91

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

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