

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097172.D
 Acq On : 3 Aug 2018 9:14
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.485

Quant Time: Aug 03 11:26:12 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 : Fri Aug 03 11:26:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1147	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5535	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3334	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9269	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9453	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8538	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3629	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	11083	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	5213	0.389	ng/ul#	92
5) 2-Methylnaphthalene	11.80	142	3569	0.403	ng/ul	98
7) Acenaphthylene	13.73	152	5777	0.401	ng/ul	96
8) Acenaphthene	14.08	153	4115	0.395	ng/ul#	93
9) Fluorene	15.08	166	4794	0.400	ng/ul	93
11) Pentachlorophenol	16.44	266	440	0.331	ng/ul	97
12) Phenanthrene	16.81	178	9112	0.386	ng/ul#	89
13) Anthracene	16.90	178	8225	0.373	ng/ul#	92
15) Fluoranthene	18.84	202	11046	0.354	ng/ul	84
17) Pyrene	19.20	202	10623	0.419	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	10094	0.374	ng/ul#	84
19) Chrysene	21.01	228	10874	0.391	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	9423	0.405	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	10059	0.401	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	9114	0.391	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	11312	0.388	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.53	278	9296	0.390	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	9922	0.394	ng/ul#	92

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

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