

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080918\
 Data File : BE097253.D
 Acq On : 9 Aug 2018 16:29
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.496

Quant Time: Aug 09 17:27:17 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 09 16:11:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1148	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6097	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3418	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	7689	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	7479	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7751	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	3878	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	8476	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	5721	0.387	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	3853	0.395	ng/ul	98
7) Acenaphthylene	13.73	152	5885	0.398	ng/ul	97
8) Acenaphthene	14.08	153	4176	0.391	ng/ul	94
9) Fluorene	15.08	166	4604	0.375	ng/ul	93
11) Pentachlorophenol	16.44	266	281	0.255	ng/ul	92
12) Phenanthrene	16.81	178	7573	0.386	ng/ul#	89
13) Anthracene	16.90	178	7047	0.385	ng/ul#	92
15) Fluoranthene	18.84	202	8437	0.326	ng/ul	84
17) Pyrene	19.20	202	8333	0.416	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	8320	0.389	ng/ul#	86
19) Chrysene	21.01	228	8486	0.385	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	7992	0.378	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	8911	0.391	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	8049	0.381	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.51	276	10438	0.395	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.53	278	8718	0.403	ng/ul#	89
26) Benzo(g,h,i)perylene	26.21	276	8762	0.383	ng/ul#	93

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

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