

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097192.D
 Acq On : 3 Aug 2018 23:01
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.487

Quant Time: Aug 04 00:35:08 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 23:39:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1405	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7630	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4435	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	10486	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	11578	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	12112	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	5005	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12762	0.35	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	7111	0.385	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4922	0.403	ng/ul	98
7) Acenaphthylene	13.74	152	7648	0.399	ng/ul	96
8) Acenaphthene	14.08	153	5438	0.392	ng/ul	96
9) Fluorene	15.08	166	6167	0.387	ng/ul#	91
11) Pentachlorophenol	16.44	266	476	0.317	ng/ul	97
12) Phenanthrene	16.81	178	10290	0.385	ng/ul#	88
13) Anthracene	16.90	178	10130	0.406	ng/ul#	91
15) Fluoranthene	18.84	202	12339	0.349	ng/ul	84
17) Pyrene	19.21	202	12769	0.412	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	13494	0.408	ng/ul#	87
19) Chrysene	21.01	228	13423	0.394	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	13365	0.405	ng/ul	89
22) Benzo(k)fluoranthene	22.58	252	13289	0.373	ng/ul	93
23) Benzo(a)pyrene	23.12	252	12970	0.393	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.51	276	16291	0.394	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.53	278	13552	0.401	ng/ul	93
26) Benzo(g,h,i)perylene	26.21	276	13653	0.382	ng/ul#	92

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

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