

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE082418\
 Data File : BE097372.D
 Acq On : 24 Aug 2018 17:19
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.412

Quant Time: Aug 24 22:51:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 24 22:49:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	758	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3599	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2009	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	5840	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	7535	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	6864	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2308	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	7832	0.50	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.17	128	3593	0.409	ng/ul#	95
5) 2-Methylnaphthalene	11.80	142	2287	0.404	ng/ul	98
7) Acenaphthylene	13.73	152	3053	0.337	ng/ul	99
8) Acenaphthene	14.07	153	2561	0.392	ng/ul	96
9) Fluorene	15.07	166	2947	0.415	ng/ul	94
11) Pentachlorophenol	16.44	266	222	0.586	ng/ul	97
12) Phenanthrene	16.80	178	6057	0.401	ng/ul#	89
13) Anthracene	16.90	178	5252	0.415	ng/ul#	91
15) Fluoranthene	18.84	202	7857	0.490	ng/ul	84
17) Pyrene	19.20	202	7775	0.341	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	8006	0.429	ng/ul#	86
19) Chrysene	21.01	228	9016	0.404	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	7724	0.416	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	8566	0.371	ng/ul#	86
23) Benzo(a)pyrene	23.11	252	7210	0.420	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.50	276	9123	0.427	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.52	278	7678	0.445	ng/ul#	89
26) Benzo(g,h,i)perylene	26.19	276	7912	0.463	ng/ul#	93

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

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