

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096112.D
 Acq On : 24 Apr 2018 7:23
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.450

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:11:34 PM

Quant Time: Apr 24 17:30:48 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 24 06:59:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	2017	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3122	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6846	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	8134	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6714m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3920	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9563	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	5868	0.370	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	3989	0.371	ng/ul	99
7) Acenaphthylene	14.70	152	6035	0.373	ng/ul	96
8) Acenaphthene	15.03	153	4313	0.371	ng/ul	93
9) Fluorene	15.99	166	4643	0.354	ng/ul	90
11) Pentachlorophenol	17.34	266	307	0.212	ng/ul	93
12) Phenanthrene	17.72	178	7428	0.368	ng/ul#	89
13) Anthracene	17.80	178	7446	0.371	ng/ul#	90
15) Fluoranthene	19.68	202	9365	0.341	ng/ul	84
17) Pyrene	20.04	202	9562	0.346	ng/ul#	79
18) Benzo(a)anthracene	21.74	228	10432	0.317	ng/ul#	84
19) Chrysene	21.80	228	8167	0.345	ng/ul	95
21) Benzo(b)fluoranthene	23.57	252	8052m	0.371	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	7804m	0.361	ng/ul	
23) Benzo(a)pyrene	24.26	252	7651m	0.365	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.17	276	8818m	0.378	ng/ul	
25) Dibenzo(a,h)anthracene	27.18	278	7126m	0.380	ng/ul	
26) Benzo(g,h,i)perylene	28.04	276	7221m	0.371	ng/ul	

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

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