

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081717\
 Data File : BE093869.D
 Acq On : 17 Aug 2017 17:29
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.413

Manual Integrations
 APPROVED

Sohil
 8/18/2017 7:08:42 PM

Quant Time: Aug 17 20:06:20 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 19:06:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1903	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	9835	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5411	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12069	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13280	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11892m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6083	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	13870	0.36	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.75	128	9437	0.394	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6406	0.367	ng/ul	98
7) Acenaphthylene	14.24	152	9502	0.420	ng/ul	97
8) Acenaphthene	14.59	153	6866	0.386	ng/ul	99
9) Fluorene	15.57	166	7724	0.343	ng/ul#	93
11) Pentachlorophenol	16.93	266	786	0.395	ng/ul	96
12) Phenanthrene	17.29	178	12362m	0.376	ng/ul	
13) Anthracene	17.39	178	12348	0.386	ng/ul	93
15) Fluoranthene	19.31	202	14700	0.358	ng/ul	85
17) Pyrene	19.67	202	15485	0.399	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	14771	0.384	ng/ul#	86
19) Chrysene	21.45	228	14652	0.397	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	13541m	0.368	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	14581m	0.402	ng/ul	
23) Benzo(a)pyrene	23.81	252	13538m	0.387	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	13866m	0.359	ng/ul	
25) Dibenzo(a,h)anthracene	26.59	278	11192m	0.347	ng/ul	
26) Benzo(g,h,i)perylene	27.38	276	11555m	0.355	ng/ul	

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

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