

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096089.D
 Acq On : 23 Apr 2018 17:37
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.448

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:10:03 PM

Quant Time: Apr 24 06:57:05 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 05:21:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1866	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5079	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2807	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	6299	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7799	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	6332m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3565	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	8891	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	5257	0.371	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	3611	0.376	ng/ul	99
7) Acenaphthylene	14.69	152	5400	0.371	ng/ul	97
8) Acenaphthene	15.03	153	3856	0.369	ng/ul	93
9) Fluorene	16.00	166	4279	0.363	ng/ul#	91
11) Pentachlorophenol	17.33	266	322	0.242	ng/ul	96
12) Phenanthrene	17.71	178	6912	0.372	ng/ul#	88
13) Anthracene	17.80	178	6719	0.364	ng/ul#	91
15) Fluoranthene	19.68	202	8658	0.343	ng/ul	85
17) Pyrene	20.04	202	9461	0.357	ng/ul#	79
18) Benzo(a)anthracene	21.74	228	10658	0.338	ng/ul#	85
19) Chrysene	21.80	228	7717	0.340	ng/ul	96
21) Benzo(b)fluoranthene	23.57	252	7693m	0.375	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	7110m	0.348	ng/ul	
23) Benzo(a)pyrene	24.26	252	7169m	0.363	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.17	276	8051m	0.366	ng/ul	
25) Dibenzo(a,h)anthracene	27.18	278	6653m	0.376	ng/ul	
26) Benzo(g,h,i)perylene	28.03	276	6808m	0.371	ng/ul	

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

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