

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094632.D
 Acq On : 28 Oct 2017 20:49
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.464

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:18 PM

Quant Time: Oct 30 04:27:54 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 02:53:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1123	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6614	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3608	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8012	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8090	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7897	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4170	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9676	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	6203	0.384	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	4379	0.387	ng/ul	98
7) Acenaphthylene	14.24	152	6661	0.408	ng/ul	97
8) Acenaphthene	14.58	153	4643	0.396	ng/ul	97
9) Fluorene	15.57	166	5197	0.395	ng/ul	91
11) Pentachlorophenol	16.98	266	223m	0.557	ng/ul	
12) Phenanthrene	17.31	178	8134	0.384	ng/ul	95
13) Anthracene	17.40	178	8565m	0.410	ng/ul	
15) Fluoranthene	19.32	202	9916	0.405	ng/ul	84
17) Pyrene	19.68	202	10469	0.419	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	9839	0.425	ng/ul#	89
19) Chrysene	21.47	228	9344	0.391	ng/ul	96
21) Benzo(b)fluoranthene	23.19	252	8890	0.400	ng/ul	94
22) Benzo(k)fluoranthene	23.24	252	9566	0.402	ng/ul#	94
23) Benzo(a)pyrene	23.86	252	8953	0.403	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	26.67	276	8584	0.374	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.67	278	7570	0.391	ng/ul#	92
26) Benzo(g,h,i)perylene	27.51	276	6305	0.342	ng/ul#	86

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

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