

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081417\
 Data File : BE093792.D
 Acq On : 14 Aug 2017 18:13
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.404

Quant Time: Aug 14 18:47:29 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 12 02:01:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2133	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11180	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6457	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14857	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16300	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	14538	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7036	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	17493	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	10788	0.397	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7574	0.382	ng/ul	99
7) Acenaphthylene	14.24	152	11591	0.429	ng/ul	96
8) Acenaphthene	14.59	153	8378	0.394	ng/ul	98
9) Fluorene	15.57	166	9687	0.361	ng/ul#	90
11) Pentachlorophenol	16.91	266	1291	0.527	ng/ul	96
12) Phenanthrene	17.29	178	15641	0.387	ng/ul#	89
13) Anthracene	17.38	178	15969	0.405	ng/ul#	88
15) Fluoranthene	19.30	202	18220	0.360	ng/ul	85
17) Pyrene	19.66	202	19231	0.404	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	18562	0.393	ng/ul#	87
19) Chrysene	21.45	228	17678	0.390	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	16584	0.369	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	17513	0.395	ng/ul	92
23) Benzo(a)pyrene	23.80	252	16872	0.394	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.56	276	17509	0.371	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.58	278	14802	0.375	ng/ul	92
26) Benzo(g,h,i)perylene	27.36	276	14104	0.354	ng/ul	99

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

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