

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093534.D
 Acq On : 24 Jul 2017 12:32
 Operator : SJ/JU
 Sample : SSTD3.234
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD3.234

Quant Time: Jul 24 13:08:03 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 11:37:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2926	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	15065	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	9820	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	25381	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	30857	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	26336	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	81015	3.67	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	262462	3.80	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	118538	3.171	ng/ul#	91
5) 2-Methylnaphthalene	12.37	142	89010	3.482	ng/ul	98
7) Acenaphthylene	14.25	152	144323	3.425	ng/ul	96
8) Acenaphthene	14.60	153	103269	3.038	ng/ul	97
9) Fluorene	15.58	166	129119	3.239	ng/ul#	90
11) Pentachlorophenol	16.93	266	15085	4.006	ng/ul	96
12) Phenanthrene	17.29	178	220527	3.158	ng/ul#	82
13) Anthracene	17.39	178	225011	3.701	ng/ul#	91
15) Fluoranthene	19.31	202	281805	3.291	ng/ul	83
17) Pyrene	19.67	202	289128	3.177	ng/ul#	79
18) Benzo(a)anthracene	21.40	228	280599	3.409	ng/ul#	86
19) Chrysene	21.45	228	267277	3.130	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	262918	3.290	ng/ul	83
22) Benzo(k)fluoranthene	23.20	252	261699	3.083	ng/ul#	84
23) Benzo(a)pyrene	23.81	252	252369	3.340	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	26.57	276	279642	3.196	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.58	278	235684	3.198	ng/ul#	87
26) Benzo(g,h,i)perylene	27.38	276	231469	3.085	ng/ul	94

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

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