

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092017\
 Data File : BE094029.D
 Acq On : 20 Sep 2017 12:02
 Operator : SJ/JU
 Sample : SSTD3.207
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD3.207

Quant Time: Sep 20 13:03:03 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 20 11:55:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1437	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7051	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4185	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.26	188	10980	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	14594	0.40	ng/ul	-0.01
20) Perylene-d12	23.92	264	12263	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	36898	3.23	ng/ul	-0.03
14) Fluoranthene-d10	19.27	212	124021	3.58	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	54779	3.194	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	39132	3.128	ng/ul	99
7) Acenaphthylene	14.24	152	59408	3.394	ng/ul	95
8) Acenaphthene	14.58	153	44294	3.216	ng/ul	98
9) Fluorene	15.57	166	54306	3.119	ng/ul#	92
11) Pentachlorophenol	16.93	266	4616	2.549	ng/ul	97
12) Phenanthrene	17.29	178	97779	3.270	ng/ul#	90
13) Anthracene	17.38	178	98763	3.391	ng/ul#	90
15) Fluoranthene	19.30	202	130609	3.493	ng/ul	84
17) Pyrene	19.66	202	134495	3.157	ng/ul#	83
18) Benzo(a)anthracene	21.40	228	135381	3.200	ng/ul#	86
19) Chrysene	21.45	228	137343	3.384	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	120703	3.182	ng/ul	83
22) Benzo(k)fluoranthene	23.20	252	130680	3.493	ng/ul#	85
23) Benzo(a)pyrene	23.81	252	118427	3.282	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.57	276	126546	3.177	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.60	278	106672	3.205	ng/ul#	88
26) Benzo(g,h,i)perylene	27.39	276	105476	3.142	ng/ul	95

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

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