

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094731.D
 Acq On : 4 Nov 2017 5:10
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.474

Quant Time: Nov 04 05:41:38 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 02:38:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1091	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6752	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3650	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7845	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8095	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7694	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3979	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	8992	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	6007	0.364	ng/ul#	94
5) 2-Methylnaphthalene	12.37	142	4094	0.354	ng/ul	99
7) Acenaphthylene	14.24	152	6078	0.368	ng/ul	98
8) Acenaphthene	14.58	153	4370	0.368	ng/ul	98
9) Fluorene	15.58	166	4839	0.363	ng/ul	91
11) Pentachlorophenol	17.01	266	176	0.449	ng/ul	96
12) Phenanthrene	17.31	178	7825	0.378	ng/ul#	93
13) Anthracene	17.39	178	7991	0.391	ng/ul#	92
15) Fluoranthene	19.32	202	9411	0.393	ng/ul	84
17) Pyrene	19.68	202	9949	0.398	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	9018	0.390	ng/ul#	87
19) Chrysene	21.47	228	8841	0.370	ng/ul	97
21) Benzo(b)fluoranthene	23.19	252	8181	0.378	ng/ul	92
22) Benzo(k)fluoranthene	23.23	252	8793	0.379	ng/ul#	93
23) Benzo(a)pyrene	23.86	252	8243	0.381	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.67	276	7505	0.335	ng/ul#	80
25) Dibenzo(a,h)anthracene	26.67	278	6431	0.341	ng/ul#	93
26) Benzo(g,h,i)perylene	27.51	276	5689	0.316	ng/ul#	89

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

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