

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094761.D
 Acq On : 6 Nov 2017 19:50
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.478

Quant Time: Nov 07 00:15:08 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 00:12:32 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1121	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	6829	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3713	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	7972	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8011	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	7702	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4060	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9174	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	5919	0.354	ng/ul#	94
5) 2-Methylnaphthalene	12.37	142	4089	0.350	ng/ul	99
7) Acenaphthylene	14.25	152	6136	0.365	ng/ul	97
8) Acenaphthene	14.59	153	4359	0.361	ng/ul	99
9) Fluorene	15.58	166	4807	0.355	ng/ul	94
11) Pentachlorophenol	17.05	266	140	0.352	ng/ul#	84
12) Phenanthrene	17.31	178	7467	0.355	ng/ul#	92
13) Anthracene	17.40	178	7893	0.380	ng/ul#	91
15) Fluoranthene	19.32	202	9311	0.382	ng/ul	84
17) Pyrene	19.68	202	9788	0.395	ng/ul#	81
18) Benzo(a)anthracene	21.42	228	8948	0.391	ng/ul#	89
19) Chrysene	21.47	228	8768	0.370	ng/ul	94
21) Benzo(b)fluoranthene	23.19	252	8354	0.386	ng/ul	93
22) Benzo(k)fluoranthene	23.24	252	8620	0.371	ng/ul#	94
23) Benzo(a)pyrene	23.86	252	8260	0.381	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.70	276	7486	0.334	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.68	278	6716	0.356	ng/ul#	88
26) Benzo(g,h,i)perylene	27.54	276	5624	0.313	ng/ul#	80

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

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