

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094665.D
 Acq On : 31 Oct 2017 19:29
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.469

Quant Time: Nov 01 01:47:39 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 01:43:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1249	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7573	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.52	164	4177	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8864	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9354	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8870	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4819	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	10849	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	7205	0.389	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	4991	0.385	ng/ul	99
7) Acenaphthylene	14.24	152	7494	0.397	ng/ul	97
8) Acenaphthene	14.58	153	5357	0.395	ng/ul	97
9) Fluorene	15.58	166	5923	0.388	ng/ul	94
11) Pentachlorophenol	17.03	266	116	0.262	ng/ul#	84
12) Phenanthrene	17.31	178	9329	0.398	ng/ul#	94
13) Anthracene	17.40	178	9220	0.399	ng/ul#	92
15) Fluoranthene	19.32	202	10941	0.404	ng/ul	85
17) Pyrene	19.68	202	11768	0.407	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	10834	0.405	ng/ul#	87
19) Chrysene	21.47	228	10938	0.396	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	9653	0.387	ng/ul	91
22) Benzo(k)fluoranthene	23.24	252	10993	0.411	ng/ul	92
23) Benzo(a)pyrene	23.86	252	9914	0.398	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.68	276	8731	0.338	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.67	278	7716	0.355	ng/ul	94
26) Benzo(g,h,i)perylene	27.53	276	6323	0.305	ng/ul	93

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

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