

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094612.D
 Acq On : 28 Oct 2017 00:59
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.462

Quant Time: Oct 30 01:13:04 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 01:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1100	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6371	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3451	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7561	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	7731	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7512	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4027	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9196	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	6015	0.386	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	4212	0.387	ng/ul	97
7) Acenaphthylene	14.25	152	6421	0.411	ng/ul	97
8) Acenaphthene	14.58	153	4476	0.399	ng/ul	96
9) Fluorene	15.57	166	5080	0.403	ng/ul	90
11) Pentachlorophenol	17.02	266	159	0.421	ng/ul	94
12) Phenanthrene	17.31	178	7800	0.391	ng/ul#	93
13) Anthracene	17.40	178	8034	0.408	ng/ul#	92
15) Fluoranthene	19.32	202	9507	0.412	ng/ul	84
17) Pyrene	19.68	202	10129	0.424	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	9394	0.425	ng/ul#	88
19) Chrysene	21.47	228	9118	0.399	ng/ul	95
21) Benzo(b)fluoranthene	23.19	252	8592	0.407	ng/ul	92
22) Benzo(k)fluoranthene	23.24	252	9030	0.399	ng/ul#	91
23) Benzo(a)pyrene	23.86	252	8596	0.407	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.67	276	8494	0.389	ng/ul#	82
25) Dibenzo(a,h)anthracene	26.67	278	7407	0.402	ng/ul#	90
26) Benzo(g,h,i)perylene	27.52	276	6335	0.361	ng/ul#	83

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

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