

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097438.D
 Acq On : 8 Sep 2018 3:07
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.428

Quant Time: Sep 08 04:17:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 08 02:48:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	864	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4599	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3108	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8320	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9326	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9584	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	3244	0.45	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	10769	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.17	128	4522	0.403	ng/ul#	94
5) 2-Methylnaphthalene	11.79	142	3259	0.460	ng/ul	98
7) Acenaphthylene	13.72	152	5076	0.448	ng/ul	97
8) Acenaphthene	14.07	153	3987	0.409	ng/ul	94
9) Fluorene	15.07	166	4707	0.422	ng/ul#	92
11) Pentachlorophenol	16.43	266	486	0.593	ng/ul	98
12) Phenanthrene	16.80	178	8659	0.404	ng/ul#	89
13) Anthracene	16.90	178	7998	0.444	ng/ul#	92
15) Fluoranthene	18.83	202	10584	0.385	ng/ul	84
17) Pyrene	19.20	202	10565	0.421	ng/ul#	81
18) Benzo(a)anthracene	20.95	228	10784	0.450	ng/ul#	85
19) Chrysene	21.01	228	10884	0.394	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	10674	0.416	ng/ul	84
22) Benzo(k)fluoranthene	22.57	252	11038	0.386	ng/ul#	92
23) Benzo(a)pyrene	23.11	252	10478	0.441	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.49	276	13130	0.427	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.51	278	11108	0.426	ng/ul#	90
26) Benzo(g,h,i)perylene	26.19	276	10965	0.418	ng/ul#	92

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

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