

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098572.D
 Acq On : 28 Dec 2018 13:45
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.409

Quant Time: Dec 28 16:08:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 13:09:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	864	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	4157	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2643	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6775	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7763	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7727	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	2656	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	8178	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	4455	0.393	ng/ul	100
5) 2-Methylnaphthalene	12.28	142	2831	0.394	ng/ul	100
7) Acenaphthylene	14.19	152	4386	0.387	ng/ul	99
8) Acenaphthene	14.53	153	3570	0.401	ng/ul	97
9) Fluorene	15.53	166	4158	0.393	ng/ul	99
11) Pentachlorophenol	16.89	266	303	0.301	ng/ul	95
12) Phenanthrene	17.27	178	7032	0.386	ng/ul	98
13) Anthracene	17.37	178	5855	0.384	ng/ul	98
15) Fluoranthene	19.29	202	8964	0.381	ng/ul	93
17) Pyrene	19.65	202	10231	0.412	ng/ul#	95
18) Benzo(a)anthracene	21.39	228	8651	0.382	ng/ul	99
19) Chrysene	21.45	228	11189	0.398	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	9223	0.390	ng/ul	99
22) Benzo(k)fluoranthene	23.20	252	10066	0.388	ng/ul	99
23) Benzo(a)pyrene	23.81	252	8857	0.384	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.57	276	10356	0.391	ng/ul#	90
25) Dibenzo(a,h)anthracene	26.60	278	8601	0.398	ng/ul	98
26) Benzo(g,h,i)perylene	27.38	276	9070	0.395	ng/ul	97

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

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