

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025495.D
 Acq On : 11 Mar 2020 17:29
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.436

Quant Time: Mar 11 18:34:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 11 07:15:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3673	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	14743	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9449	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	18994	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15220	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	16020	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	8882	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	16918	0.32	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.48	128	14709	0.337	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	10325	0.331	ng/ul	100
7) Acenaphthylene	14.01	152	11945	0.309	ng/ul	99
8) Acenaphthene	14.35	153	11026	0.326	ng/ul	100
9) Fluorene	15.34	166	12746	0.309	ng/ul	99
11) Pentachlorophenol	16.69	266	1463	0.328	ng/ul	98
12) Phenanthrene	17.07	178	19501	0.324	ng/ul	99
13) Anthracene	17.16	178	16916	0.323	ng/ul	99
15) Fluoranthene	19.09	202	20770	0.318	ng/ul	98
17) Pyrene	19.45	202	20439	0.299	ng/ul	98
18) Benzo(a)anthracene	21.20	228	18109	0.333	ng/ul	99
19) Chrysene	21.26	228	20159	0.330	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	19548	0.308	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	21052	0.317	ng/ul	99
23) Benzo(a)pyrene	23.31	252	17362	0.330	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	23506	0.329	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.59	278	19644	0.327	ng/ul	98
26) Benzo(g,h,i)perylene	26.24	276	20059	0.324	ng/ul	98

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

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