

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025477.D
 Acq On : 11 Mar 2020 06:34
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 11 07:14:45 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 06:51:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3807	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15778	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	10959	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	24076	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	16619	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	16238	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	9793	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	20880	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	15833	0.339	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	11401	0.342	ng/ul	100
7) Acenaphthylene	14.01	152	13936	0.311	ng/ul	99
8) Acenaphthene	14.35	153	13081	0.333	ng/ul	99
9) Fluorene	15.34	166	15629	0.327	ng/ul	98
11) Pentachlorophenol	16.69	266	1798	0.318	ng/ul	98
12) Phenanthrene	17.07	178	24872	0.326	ng/ul	100
13) Anthracene	17.16	178	21461	0.323	ng/ul	99
15) Fluoranthene	19.09	202	25825	0.312	ng/ul	98
17) Pyrene	19.45	202	25111	0.337	ng/ul	99
18) Benzo(a)anthracene	21.20	228	19746	0.332	ng/ul	99
19) Chrysene	21.26	228	21860	0.328	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	20393	0.317	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	21189m	0.315	ng/ul	
23) Benzo(a)pyrene	23.31	252	17568	0.329	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.58	276	24190	0.334	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.59	278	19916	0.327	ng/ul	98
26) Benzo(g,h,i)perylene	26.23	276	19932	0.317	ng/ul	97

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

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