

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025412.D
 Acq On : 09 Mar 2020 11:46
 Operator : CG/JU
 Sample : SST0.876
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.876

Quant Time: Mar 09 13:20:22 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 : Mon Mar 09 12:03:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2122	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	8213	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	5570	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	12222	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8983	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	8229	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	12446	0.81	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	27470	0.69	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.48	128	20080	0.812	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	14537	0.807	ng/ul	100
7) Acenaphthylene	14.01	152	18629	0.714	ng/ul	99
8) Acenaphthene	14.35	153	16731	0.827	ng/ul	99
9) Fluorene	15.34	166	20559	0.811	ng/ul	99
11) Pentachlorophenol	16.69	266	2292	0.716	ng/ul	99
12) Phenanthrene	17.07	178	32462	0.824	ng/ul	100
13) Anthracene	17.16	178	28725	0.821	ng/ul	99
15) Fluoranthene	19.09	202	34437	0.698	ng/ul	99
17) Pyrene	19.46	202	33041	0.970	ng/ul	98
18) Benzo(a)anthracene	21.20	228	26726	0.809	ng/ul	99
19) Chrysene	21.26	228	29788	0.821	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	27378	0.832	ng/ul	96
22) Benzo(k)fluoranthene	22.80	252	28387	0.822	ng/ul	97
23) Benzo(a)pyrene	23.31	252	22961	0.831	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.58	276	31608	0.917	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.59	278	26764	0.937	ng/ul	98
26) Benzo(g,h,i)perylene	26.24	276	27457	0.938	ng/ul	99

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

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