

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025410.D
 Acq On : 09 Mar 2020 10:34
 Operator : CG/JU
 Sample : SSTDO.280
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDO.280

Quant Time: Mar 09 12:07:18 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	7.66	152	2335	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	9170	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	6161	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	14055	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	9923	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	8989	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	3497	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	8115	0.18	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.48	128	5765	0.209	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	4031	0.201	ng/ul	99
7) Acenaphthylene	14.01	152	5379	0.186	ng/ul	99
8) Acenaphthene	14.35	153	4650	0.208	ng/ul	98
9) Fluorene	15.34	166	5664	0.202	ng/ul	99
11) Pentachlorophenol	16.70	266	583	0.158	ng/ul	96
12) Phenanthrene	17.08	178	9321	0.206	ng/ul	98
13) Anthracene	17.17	178	7942	0.197	ng/ul	98
15) Fluoranthene	19.09	202	9984	0.176	ng/ul	99
17) Pyrene	19.45	202	9699	0.258	ng/ul	99
18) Benzo(a)anthracene	21.21	228	7407	0.203	ng/ul	99
19) Chrysene	21.26	228	8415	0.210	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	7325	0.204	ng/ul	93
22) Benzo(k)fluoranthene	22.80	252	7800	0.207	ng/ul#	92
23) Benzo(a)pyrene	23.31	252	6053	0.200	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.58	276	8221	0.218	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.60	278	6839	0.219	ng/ul	95
26) Benzo(g,h,i)perylene	26.24	276	7161	0.224	ng/ul	97

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

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