

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031320\
 Data File : BM025542.D
 Acq On : 13 Mar 2020 15:42
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.441

Quant Time: Mar 13 16:54:36 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 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3146	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	12724	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	8688	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	18108	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12075	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	12438	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	9017	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	18010	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	14342	0.381	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	10407	0.387	ng/ul	100
7) Acenaphthylene	14.00	152	12595	0.355	ng/ul#	86
8) Acenaphthene	14.35	153	11893	0.382	ng/ul	99
9) Fluorene	15.33	166	14330	0.378	ng/ul	99
11) Pentachlorophenol	16.69	266	1560	0.367	ng/ul	98
12) Phenanthrene	17.07	178	22128	0.386	ng/ul	100
13) Anthracene	17.16	178	18825	0.377	ng/ul	100
15) Fluoranthene	19.09	202	22028	0.354	ng/ul	98
17) Pyrene	19.45	202	21482	0.396	ng/ul	98
18) Benzo(a)anthracene	21.20	228	17074	0.395	ng/ul	99
19) Chrysene	21.25	228	19288	0.398	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	18228	0.370	ng/ul	98
22) Benzo(k)fluoranthene	22.80	252	19065	0.369	ng/ul	99
23) Benzo(a)pyrene	23.31	252	15807	0.386	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.57	276	20340	0.367	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.59	278	16631	0.356	ng/ul	99
26) Benzo(g,h,i)perylene	26.23	276	17998	0.374	ng/ul	98

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

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