

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025031.D
 Acq On : 24 Feb 2020 20:47
 Operator : CG/JU
 Sample : SST0.880
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.880

Quant Time: Feb 25 00:51:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 00:46:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1992	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	8068	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	5604	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	13837	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15195	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	16433	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	11756	0.89	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	37091	0.98	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	18953	0.824	ng/ul#	89
5) 2-Methylnaphthalene	12.14	142	13977	0.887	ng/ul	99
7) Acenaphthylene	14.05	152	24760	0.954	ng/ul	95
8) Acenaphthene	14.39	153	16668	0.768	ng/ul	98
9) Fluorene	15.38	166	20882	0.839	ng/ul	98
11) Pentachlorophenol	16.72	266	3324	1.184	ng/ul	97
12) Phenanthrene	17.11	178	35531	0.806	ng/ul	97
13) Anthracene	17.20	178	36325	0.960	ng/ul	97
15) Fluoranthene	19.13	202	47633	1.001	ng/ul	98
17) Pyrene	19.49	202	48137	0.699	ng/ul	97
18) Benzo(a)anthracene	21.24	228	48630	0.836	ng/ul	98
19) Chrysene	21.29	228	47940	0.784	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	49464	0.693	ng/ul	93
22) Benzo(k)fluoranthene	22.84	252	49777	0.683	ng/ul#	95
23) Benzo(a)pyrene	23.36	252	45821	0.733	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.65	276	57764	0.758	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	48207	0.783	ng/ul	95
26) Benzo(g,h,i)perylene	26.32	276	48436	0.798	ng/ul#	94

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

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