

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031320\
 Data File : BM025548.D
 Acq On : 13 Mar 2020 19:24
 Operator : CG/JU
 Sample : L1786-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC6

Quant Time: Mar 16 07:39:34 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 16 06:44:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	5	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	77	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	280	0.40	ng/ul	0.00
10) Phenanthrene-d10	0.00	188	0	0.00	ng/ul	-17.03
16) Chrysene-d12	21.21	240	392	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	14	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2037	14.67	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4718	0.00	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.47	128	60197	264.090	ng/ul	98
5) 2-Methylnaphthalene	12.09	142	13485	82.772	ng/ul	99
7) Acenaphthylene	14.00	152	28292	24.731	ng/ul#	86
8) Acenaphthene	14.35	153	2913	2.903	ng/ul	98
9) Fluorene	15.33	166	25583	20.923	ng/ul	99
17) Pyrene	19.45	202	221757	126.037	ng/ul	100
18) Benzo(a)anthracene	21.20	228	188043	134.171	ng/ul	98
19) Chrysene	21.25	228	154092	97.909	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	285617	5152.062	ng/ul	96
22) Benzo(k)fluoranthene	22.75	252	285617	4917.228	ng/ul	96
23) Benzo(a)pyrene	23.30	252	107726	2340.037	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.56	276	72128	1154.839	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.57	278	24565	467.535	ng/ul	97
26) Benzo(g,h,i)perylene	26.23	276	39514	729.338	ng/ul	98

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

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