

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101220\
 Data File : BM027654.D
 Acq On : 12 Oct 2020 12:03
 Operator : JU/CG
 Sample : PB132211BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS211

Quant Time: Oct 12 15:10:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM100820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 10:50:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	610	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	1862	0.40	ng/ul	0.00
7) Acenaphthene-d10	14.41	164	1341	0.40	ng/ul	0.00
11) Phenanthrene-d10	17.16	188	3572	0.40	ng/ul	0.00
17) Chrysene-d12	21.34	240	3898	0.40	ng/ul	0.00
21) Perylene-d12	23.59	264	4128	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.16	152	1213	0.35	ng/ul	0.00
15) Fluoranthene-d10	19.18	212	4303	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.61	128	1713	0.323	ng/ul	96
5) 2-Methylnaphthalene	12.23	142	1262	0.328	ng/ul	98
6) 1-Methylnaphthalene	12.45	142	1451	0.362	ng/ul	100
8) Acenaphthylene	14.13	152	1911	0.328	ng/ul	99
9) Acenaphthene	14.48	153	1521	0.330	ng/ul	97
10) Fluorene	15.46	166	1989	0.339	ng/ul	99
12) Pentachlorophenol	16.80	266	998	0.688	ng/ul	98
13) Phenanthrene	17.20	178	3615	0.334	ng/ul	97
14) Anthracene	17.29	178	3387	0.335	ng/ul	99
16) Fluoranthene	19.21	202	5051	0.370	ng/ul	97
18) Pyrene	19.57	202	5154	0.382	ng/ul#	97
19) Benzo(a)anthracene	21.32	228	5421	0.359	ng/ul	98
20) Chrysene	21.37	228	5702	0.356	ng/ul	99
22) Benzo(b)fluoranthene	22.91	252	6198	0.354	ng/ul	96
23) Benzo(k)fluoranthene	22.96	252	6083	0.346	ng/ul	97
24) Benzo(a)pyrene	23.49	252	5535	0.360	ng/ul#	94
25) Indeno(1,2,3-cd)pyrene	25.85	276	7211	0.349	ng/ul#	98
26) Dibenzo(a,h)anthracene	25.87	278	6012	0.352	ng/ul	97
27) Benzo(a,h,i)perylene	26.55	276	6195	0.342	ng/ul	98

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

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