

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071122\
 Data File : BN020706.D
 Acq On : 11 Jul 2022 17:15
 Operator : CG/JU
 Sample : PB146063BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS063

Quant Time: Jul 11 18:40:44 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 02 00:47:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	3960	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	12557	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	6740	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	12622	0.400	ng/ul	0.00
17) Chrysene-d12	21.433	240	8260	0.400	ng/ul	-0.01
23) Perylene-d12	23.789	264	5628	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	3690	0.804	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	7517	0.365	ng/ul	0.00
18) Fluoranthene-d10	19.267	212	13146	0.483	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.256	88	9472	2.022	ng/ul#	79
5) Naphthalene	10.672	128	14382	0.371	ng/ul	99
7) 2-Methylnaphthalene	12.294	142	8670	0.334	ng/ul	99
8) 1-Methylnaphthalene	12.514	142	9222	0.375	ng/ul	100
10) Acenaphthylene	14.192	152	11552	0.362	ng/ul	99
11) Acenaphthene	14.539	153	9470	0.370	ng/ul	99
12) Fluorene	15.529	166	10244	0.338	ng/ul	100
14) Pentachlorophenol	16.892	266	1033	0.406	ng/ul	99
15) Phenanthrene	17.268	178	16421	0.374	ng/ul	99
16) Anthracene	17.365	178	12792	0.338	ng/ul	100
19) Fluoranthene	19.295	202	17081	0.453	ng/ul	99
20) Pyrene	19.657	202	17392	0.444	ng/ul	97
21) Benzo(a)anthracene	21.422	228	10176	0.343	ng/ul	100
22) Chrysene	21.474	228	13486	0.408	ng/ul	100
24) Benzo(b)fluoranthene	23.073	252	11173	0.429	ng/ul	98
25) Benzo(k)fluoranthene	23.120	252	11498	0.437	ng/ul	99
26) Benzo(a)pyrene	23.687	252	9516	0.402	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.226	276	10387	0.441	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.243	278	8227	0.432	ng/ul	98
29) Benzo(g,h,i)perylene	26.968	276	9616	0.476	ng/ul	100

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

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