

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020971.D
 Acq On : 03 Aug 2022 10:08
 Operator : CG/JU
 Sample : PB146572BS
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS572

Quant Time: Aug 03 13:41:55 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 04:19:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.809	152	1952	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	6261	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	3526	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	6949	0.400	ng/ul	0.00
17) Chrysene-d12	21.421	240	6115	0.400	ng/ul	0.00
23) Perylene-d12	23.762	264	5520	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1583	0.665	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.208	152	3252	0.352	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	6887	0.368	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	4076	1.590	ng/ul#	86
5) Naphthalene	10.651	128	5980	0.349	ng/ul	98
7) 2-Methylnaphthalene	12.279	142	3607	0.346	ng/ul	99
8) 1-Methylnaphthalene	12.499	142	3931	0.342	ng/ul	99
10) Acenaphthylene	14.179	152	5167	0.358	ng/ul	99
11) Acenaphthene	14.521	153	4271	0.355	ng/ul	99
12) Fluorene	15.511	166	4656	0.336	ng/ul	98
14) Pentachlorophenol	16.876	266	663	0.598	ng/ul	99
15) Phenanthrene	17.252	178	7590	0.322	ng/ul	100
16) Anthracene	17.349	178	6121	0.302	ng/ul	99
19) Fluoranthene	19.277	202	8467	0.308	ng/ul	98
20) Pyrene	19.644	202	8772	0.305	ng/ul	98
21) Benzo(a)anthracene	21.404	228	6502	0.293	ng/ul	99
22) Chrysene	21.456	228	8418	0.307	ng/ul	99
24) Benzo(b)fluoranthene	23.049	252	8005	0.304	ng/ul	100
25) Benzo(k)fluoranthene	23.096	252	8571	0.300	ng/ul	99
26) Benzo(a)pyrene	23.660	252	7405	0.298	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.186	276	8749	0.299	ng/ul	100
28) Dibenzo(a,h)anthracene	26.202	278	7012	0.293	ng/ul	98
29) Benzo(g,h,i)perylene	26.927	276	7954	0.299	ng/ul	96

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

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