

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062922\
 Data File : BN020579.D
 Acq On : 29 Jun 2022 20:58
 Operator : CG/JU
 Sample : PB145855BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS855

Quant Time: Jun 29 22:00:04 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 : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	3161	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.522	136	10321	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.371	164	6024	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.110	188	12367	0.400	ng/ul	0.00
17) Chrysene-d12	21.301	240	9518	0.400	ng/ul	0.00
23) Perylene-d12	23.580	264	6926	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	2479	0.677	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.117	152	5555	0.328	ng/ul	-0.01
18) Fluoranthene-d10	19.145	212	11699	0.373	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.226	88	6581	1.760	ng/ul#	80
5) Naphthalene	10.572	128	10460	0.328	ng/ul	99
7) 2-Methylnaphthalene	12.194	142	6522	0.306	ng/ul	99
8) 1-Methylnaphthalene	12.409	142	6954	0.344	ng/ul	99
10) Acenaphthylene	14.089	152	9419	0.330	ng/ul	100
11) Acenaphthene	14.431	153	7564	0.331	ng/ul	99
12) Fluorene	15.421	166	8470	0.313	ng/ul	99
14) Pentachlorophenol	16.777	266	1085	0.435	ng/ul	98
15) Phenanthrene	17.152	178	14297	0.333	ng/ul	100
16) Anthracene	17.245	178	11867	0.320	ng/ul	100
19) Fluoranthene	19.172	202	15548	0.358	ng/ul	100
20) Pyrene	19.535	202	15891	0.352	ng/ul	98
21) Benzo(a)anthracene	21.286	228	11200	0.328	ng/ul	99
22) Chrysene	21.336	228	13730	0.361	ng/ul	99
24) Benzo(b)fluoranthene	22.893	252	11543	0.360	ng/ul	98
25) Benzo(k)fluoranthene	22.937	252	11821	0.365	ng/ul	99
26) Benzo(a)pyrene	23.481	252	10414	0.358	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.917	276	10596	0.366	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.930	278	8526	0.364	ng/ul	98
29) Benzo(g,h,i)perylene	26.631	276	9560	0.384	ng/ul	99

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

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