

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020440.D
 Acq On : 25 Jun 2022 18:05
 Operator : CG/JU
 Sample : PB145678BS
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS678

Quant Time: Jun 27 03:42:53 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	7.750	152	2378	0.400	ng/u1	0.00
4) Naphthalene-d8	10.535	136	8150	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.377	164	5433	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	12417	0.400	ng/u1	0.00
17) Chrysene-d12	21.309	240	11246	0.400	ng/u1	0.00
23) Perylene-d12	23.589	264	9383	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	2023	0.734	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.130	152	5028	0.376	ng/u1	0.00
18) Fluoranthene-d10	19.151	212	13244	0.357	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.235	88	4791	1.703	ng/u1#	84
5) Naphthalene	10.584	128	8375	0.333	ng/u1	100
7) 2-Methylnaphthalene	12.201	142	5617	0.333	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	5961	0.373	ng/u1	100
10) Acenaphthylene	14.095	152	8782	0.341	ng/u1	100
11) Acenaphthene	14.442	153	6822	0.331	ng/u1	98
12) Fluorene	15.427	166	8079	0.331	ng/u1	99
14) Pentachlorophenol	16.786	266	1302	0.520	ng/u1	99
15) Phenanthrene	17.162	178	14192	0.329	ng/u1	100
16) Anthracene	17.251	178	12473	0.335	ng/u1	99
19) Fluoranthene	19.179	202	16380	0.319	ng/u1	99
20) Pyrene	19.541	202	17199	0.323	ng/u1	99
21) Benzo(a)anthracene	21.292	228	14147	0.350	ng/u1	100
22) Chrysene	21.344	228	15915	0.354	ng/u1	99
24) Benzo(b)fluoranthene	22.902	252	14819	0.341	ng/u1	99
25) Benzo(k)fluoranthene	22.946	252	15102	0.345	ng/u1	99
26) Benzo(a)pyrene	23.490	252	13963	0.354	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	25.930	276	14280	0.364	ng/u1	97
28) Dibenzo(a,h)anthracene	25.940	278	11555	0.364	ng/u1	99
29) Benzo(g,h,i)perylene	26.637	276	12310	0.365	ng/u1	99

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

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