

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020859.D
 Acq On : 26 Jul 2022 13:14
 Operator : CG/JU
 Sample : PB146369BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS369

Quant Time: Jul 26 13:52:14 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 18:06:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	2233	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	7862	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	4586	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.217	188	9497	0.400	ng/ul	0.00
17) Chrysene-d12	21.426	240	8437	0.400	ng/ul	0.00
23) Perylene-d12	23.770	264	7008	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1972	0.735	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	4816	0.403	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	11439	0.424	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4772	1.731	ng/ul#	82
5) Naphthalene	10.661	128	8681	0.352	ng/ul	100
7) 2-Methylnaphthalene	12.289	142	5344	0.343	ng/ul	99
8) 1-Methylnaphthalene	12.504	142	5766	0.382	ng/ul	100
10) Acenaphthylene	14.187	152	8049	0.353	ng/ul	100
11) Acenaphthene	14.530	153	6458	0.350	ng/ul	100
12) Fluorene	15.524	166	7209	0.356	ng/ul	99
14) Pentachlorophenol	16.884	266	1231	0.861	ng/ul	99
15) Phenanthrene	17.259	178	12034	0.346	ng/ul	99
16) Anthracene	17.356	178	9966	0.357	ng/ul	99
19) Fluoranthene	19.285	202	13941	0.353	ng/ul	99
20) Pyrene	19.648	202	14403	0.349	ng/ul	99
21) Benzo(a)anthracene	21.411	228	10402	0.373	ng/ul	100
22) Chrysene	21.461	228	12931	0.360	ng/ul	99
24) Benzo(b)fluoranthene	23.060	252	11941	0.360	ng/ul	98
25) Benzo(k)fluoranthene	23.107	252	12508	0.364	ng/ul	98
26) Benzo(a)pyrene	23.671	252	10875	0.355	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.205	276	11580	0.359	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.225	278	9072	0.369	ng/ul	99
29) Benzo(g,h,i)perylene	26.936	276	10476	0.361	ng/ul	99

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

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