

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072922\
 Data File : BN020922.D
 Acq On : 29 Jul 2022 15:02
 Operator : CG/JU
 Sample : PB146499BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS499

Quant Time: Jul 29 17:46:34 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	7.817	152	2165	0.400	ng/ul	0.00
4) Naphthalene-d8	10.611	136	7761	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.464	164	4927	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.216	188	10289	0.400	ng/ul	0.00
17) Chrysene-d12	21.424	240	8271	0.400	ng/ul	0.00
23) Perylene-d12	23.768	264	6681	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1795	0.690	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	4228	0.359	ng/ul	0.00
18) Fluoranthene-d10	19.252	212	10167	0.385	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4307	1.611	ng/ul#	82
5) Naphthalene	10.660	128	6584	0.271	ng/ul	99
7) 2-Methylnaphthalene	12.288	142	4251	0.277	ng/ul	99
8) 1-Methylnaphthalene	12.502	142	4631	0.311	ng/ul	100
10) Acenaphthylene	14.186	152	6579	0.268	ng/ul	100
11) Acenaphthene	14.524	153	5189	0.262	ng/ul	100
12) Fluorene	15.518	166	5820	0.268	ng/ul	99
14) Pentachlorophenol	16.882	266	892	0.576	ng/ul	99
15) Phenanthrene	17.258	178	9640	0.256	ng/ul	99
16) Anthracene	17.355	178	8003	0.265	ng/ul	99
19) Fluoranthene	19.284	202	11086	0.286	ng/ul	99
20) Pyrene	19.647	202	11467	0.284	ng/ul	98
21) Benzo(a)anthracene	21.412	228	7931	0.290	ng/ul	99
22) Chrysene	21.462	228	9623	0.274	ng/ul	99
24) Benzo(b)fluoranthene	23.057	252	8682	0.274	ng/ul	97
25) Benzo(k)fluoranthene	23.107	252	9069	0.277	ng/ul	98
26) Benzo(a)pyrene	23.668	252	7953	0.272	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.202	276	8100	0.264	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.222	278	6325	0.270	ng/ul	95
29) Benzo(g,h,i)perylene	26.940	276	7403	0.267	ng/ul	98

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

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