

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092122\
 Data File : BN021864.D
 Acq On : 22 Sep 2022 10:53
 Operator : CG/JU
 Sample : PB147689BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS689

Quant Time: Sep 22 12:53:22 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 00:20:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	5354	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	17758	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.707	164	9503	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.460	188	16549	0.400	ng/ul	0.00
17) Chrysene-d12	21.653	240	9635	0.400	ng/ul	# 0.00
23) Perylene-d12	24.164	264	7343	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	4140	0.629	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.467	152	9096	0.324	ng/ul	0.00
18) Fluoranthene-d10	19.487	212	12672	0.354	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.401	88	10938	1.619	ng/ul#	89
5) Naphthalene	10.927	128	16372	0.327	ng/ul	98
7) 2-Methylnaphthalene	12.539	142	10396	0.319	ng/ul	98
8) 1-Methylnaphthalene	12.759	142	10819	0.324	ng/ul	100
10) Acenaphthylene	14.430	152	13118	0.329	ng/ul	99
11) Acenaphthene	14.768	153	10850	0.321	ng/ul	100
12) Fluorene	15.758	166	11735	0.309	ng/ul	99
14) Pentachlorophenol	17.114	266	1123	0.350	ng/ul	99
15) Phenanthrene	17.502	178	17069	0.327	ng/ul	99
16) Anthracene	17.595	178	13763	0.318	ng/ul	99
19) Fluoranthene	19.520	202	15887	0.353	ng/ul	98
20) Pyrene	19.882	202	15676	0.353	ng/ul	96
21) Benzo(a)anthracene	21.638	228	11327	0.351	ng/ul	99
22) Chrysene	21.691	228	12728	0.362	ng/ul	99
24) Benzo(b)fluoranthene	23.386	252	12051	0.338	ng/ul	91
25) Benzo(k)fluoranthene	23.439	252	11057	0.320	ng/ul#	92
26) Benzo(a)pyrene	24.050	252	10568	0.392	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.807	276	11764	0.387	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.831	278	9422	0.366	ng/ul	97
29) Benzo(g,h,i)perylene	27.615	276	10655	0.381	ng/ul	96

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

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