

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092122\
 Data File : BN021901.D
 Acq On : 23 Sep 2022 14:45
 Operator : CG/JU
 Sample : PB147816BS
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS816

Quant Time: Sep 23 16:20:32 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	8.046	152	5025	0.400	ng/ul	0.00
4) Naphthalene-d8	10.872	136	16688	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.703	164	9095	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.455	188	16410	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.650	240	9788	0.400	ng/ul	# 0.00
23) Perylene-d12	24.158	264	8044	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.359	96	3287	0.532	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.461	152	7571	0.287	ng/ul	-0.01
18) Fluoranthene-d10	19.487	212	11434	0.314	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.397	88	8869	1.399	ng/ul#	87
5) Naphthalene	10.922	128	13429	0.285	ng/ul	98
7) 2-Methylnaphthalene	12.538	142	8632	0.282	ng/ul	98
8) 1-Methylnaphthalene	12.753	142	8956	0.286	ng/ul	99
10) Acenaphthylene	14.425	152	11701	0.306	ng/ul	99
11) Acenaphthene	14.768	153	9237	0.285	ng/ul	100
12) Fluorene	15.753	166	10016	0.276	ng/ul	99
14) Pentachlorophenol	17.109	266	1019	0.320	ng/ul	99
15) Phenanthrene	17.498	178	14747	0.285	ng/ul	100
16) Anthracene	17.591	178	12354	0.287	ng/ul	99
19) Fluoranthene	19.515	202	14199	0.311	ng/ul	97
20) Pyrene	19.878	202	14038	0.311	ng/ul#	94
21) Benzo(a)anthracene	21.632	228	10644	0.325	ng/ul	99
22) Chrysene	21.688	228	11412	0.319	ng/ul	99
24) Benzo(b)fluoranthene	23.386	252	10730	0.275	ng/ul	89
25) Benzo(k)fluoranthene	23.436	252	10505	0.278	ng/ul#	89
26) Benzo(a)pyrene	24.044	252	10012	0.339	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.804	276	11275	0.339	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.828	278	9190	0.326	ng/ul	96
29) Benzo(g,h,i)perylene	27.616	276	10047	0.328	ng/ul	94

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

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