

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091922\
 Data File : BN021833.D
 Acq On : 20 Sep 2022 11:14
 Operator : CG/JU
 Sample : PB147687BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS687

Quant Time: Sep 21 04:25:41 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	8.050	152	4574	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	15564	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.707	164	8595	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.460	188	15946	0.400	ng/ul	0.00
17) Chrysene-d12	21.650	240	9683	0.400	ng/ul	# 0.00
23) Perylene-d12	24.155	264	7938	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	3086	0.548	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.467	152	7251	0.294	ng/ul	0.00
18) Fluoranthene-d10	19.487	212	11404	0.317	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.401	88	8744	1.515	ng/ul#	88
5) Naphthalene	10.933	128	13128	0.299	ng/ul	98
7) 2-Methylnaphthalene	12.544	142	8487	0.298	ng/ul	98
8) 1-Methylnaphthalene	12.758	142	8853	0.303	ng/ul	99
10) Acenaphthylene	14.430	152	11388	0.315	ng/ul	99
11) Acenaphthene	14.772	153	9018	0.295	ng/ul	99
12) Fluorene	15.758	166	10039	0.293	ng/ul	99
14) Pentachlorophenol	17.113	266	937	0.303	ng/ul	99
15) Phenanthrene	17.502	178	15078	0.300	ng/ul	99
16) Anthracene	17.595	178	12662	0.303	ng/ul	99
19) Fluoranthene	19.515	202	14584	0.323	ng/ul	97
20) Pyrene	19.882	202	14394	0.323	ng/ul	96
21) Benzo(a)anthracene	21.632	228	10758	0.332	ng/ul	98
22) Chrysene	21.688	228	11394	0.322	ng/ul	99
24) Benzo(b)fluoranthene	23.383	252	11010	0.286	ng/ul	90
25) Benzo(k)fluoranthene	23.433	252	10261	0.275	ng/ul#	89
26) Benzo(a)pyrene	24.041	252	9962	0.342	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.797	276	11563	0.352	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.824	278	9425	0.338	ng/ul	96
29) Benzo(g,h,i)perylene	27.612	276	10126	0.335	ng/ul	95

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

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