

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092722\
 Data File : BN021923.D
 Acq On : 26 Sep 2022 18:17
 Operator : CG/JU
 Sample : PB147749BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS749

Quant Time: Sep 27 00:25:41 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	8.046	152	3974	0.400	ng/ul	0.00
4) Naphthalene-d8	10.872	136	13088	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.703	164	6777	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.455	188	11537	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.647	240	6775	0.400	ng/ul	# 0.00
23) Perylene-d12	24.155	264	5072	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.359	96	3317	0.678	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.467	152	7288	0.352	ng/ul	0.00
18) Fluoranthene-d10	19.487	212	9892	0.392	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.397	88	9649	1.924	ng/ul#	86
5) Naphthalene	10.927	128	13972	0.379	ng/ul	98
7) 2-Methylnaphthalene	12.538	142	8788	0.366	ng/ul	98
8) 1-Methylnaphthalene	12.753	142	9063	0.368	ng/ul	100
10) Acenaphthylene	14.425	152	11036	0.388	ng/ul	99
11) Acenaphthene	14.767	153	9192	0.381	ng/ul	100
12) Fluorene	15.757	166	9756	0.361	ng/ul	100
14) Pentachlorophenol	17.113	266	992	0.444	ng/ul	100
15) Phenanthrene	17.497	178	13727	0.377	ng/ul	99
16) Anthracene	17.590	178	10979	0.363	ng/ul	99
19) Fluoranthene	19.515	202	12995	0.411	ng/ul	97
20) Pyrene	19.877	202	12911	0.414	ng/ul	94
21) Benzo(a)anthracene	21.632	228	9220	0.406	ng/ul	99
22) Chrysene	21.685	228	10388	0.420	ng/ul	100
24) Benzo(b)fluoranthene	23.380	252	9862	0.401	ng/ul	88
25) Benzo(k)fluoranthene	23.433	252	8968	0.376	ng/ul#	85
26) Benzo(a)pyrene	24.044	252	8563	0.459	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.803	276	8914	0.425	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.824	278	6905	0.388	ng/ul#	93
29) Benzo(g,h,i)perylene	27.612	276	8550	0.443	ng/ul	94

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

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