

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080922\
 Data File : BN021095.D
 Acq On : 09 Aug 2022 21:16
 Operator : CG/JU
 Sample : PB146837BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS837

Quant Time: Aug 09 23:14:18 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 09 23:10:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	2463	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	8424	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.443	164	4831	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.193	188	9469	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	7626	0.400	ng/ul	0.00
23) Perylene-d12	23.724	264	5804	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	2807	0.891	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.185	152	6380	0.517	ng/ul	0.00
18) Fluoranthene-d10	19.231	212	13166	0.559	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.232	88	7242	2.231	ng/ul#	84
5) Naphthalene	10.635	128	12078	0.467	ng/ul	99
7) 2-Methylnaphthalene	12.262	142	7623	0.486	ng/ul	99
8) 1-Methylnaphthalene	12.477	142	7700	0.490	ng/ul	100
10) Acenaphthylene	14.160	152	11007	0.487	ng/ul	98
11) Acenaphthene	14.503	153	8459	0.453	ng/ul	99
12) Fluorene	15.493	166	9469	0.464	ng/ul	100
14) Pentachlorophenol	16.855	266	1941	1.329	ng/ul	99
15) Phenanthrene	17.235	178	15082	0.445	ng/ul	99
16) Anthracene	17.328	178	12991	0.495	ng/ul	98
19) Fluoranthene	19.259	202	16754	0.496	ng/ul	99
20) Pyrene	19.621	202	17247	0.495	ng/ul	98
21) Benzo(a)anthracene	21.377	228	12864	0.517	ng/ul	100
22) Chrysene	21.430	228	14506	0.443	ng/ul	99
24) Benzo(b)fluoranthene	23.017	252	12554	0.467	ng/ul	100
25) Benzo(k)fluoranthene	23.064	252	12790	0.451	ng/ul	99
26) Benzo(a)pyrene	23.622	252	11485	0.469	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.128	276	11432	0.391	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.152	278	9111	0.394	ng/ul	98
29) Benzo(g,h,i)perylene	26.870	276	10068	0.386	ng/ul	100

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

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