

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080922\
 Data File : BN021099.D
 Acq On : 10 Aug 2022 00:25
 Operator : CG/JU
 Sample : PB146855BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS855

Quant Time: Aug 10 01:46:28 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	2525	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	8705	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.443	164	4991	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.193	188	9482	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	7227	0.400	ng/ul	0.00
23) Perylene-d12	23.721	264	5481	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	2316	0.717	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.191	152	4931	0.387	ng/ul	0.01
18) Fluoranthene-d10	19.231	212	9735	0.436	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.232	88	5492	1.650	ng/ul#	86
5) Naphthalene	10.635	128	7853	0.294	ng/ul	99
7) 2-Methylnaphthalene	12.263	142	4868	0.300	ng/ul	99
8) 1-Methylnaphthalene	12.477	142	5258	0.324	ng/ul	100
10) Acenaphthylene	14.160	152	6930	0.297	ng/ul	99
11) Acenaphthene	14.503	153	5566	0.288	ng/ul	99
12) Fluorene	15.498	166	6091	0.289	ng/ul	100
14) Pentachlorophenol	16.859	266	821	0.562	ng/ul	98
15) Phenanthrene	17.235	178	9719	0.287	ng/ul	100
16) Anthracene	17.328	178	7978	0.304	ng/ul	99
19) Fluoranthene	19.259	202	10720	0.335	ng/ul	99
20) Pyrene	19.621	202	10989	0.333	ng/ul	97
21) Benzo(a)anthracene	21.377	228	7584	0.322	ng/ul	100
22) Chrysene	21.430	228	9110	0.293	ng/ul	99
24) Benzo(b)fluoranthene	23.014	252	7683	0.303	ng/ul	97
25) Benzo(k)fluoranthene	23.064	252	8019	0.299	ng/ul	97
26) Benzo(a)pyrene	23.622	252	7016	0.303	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.135	276	7079	0.256	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.152	278	5569	0.255	ng/ul	98
29) Benzo(g,h,i)perylene	26.870	276	6434	0.261	ng/ul	98

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

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