

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021171.D
 Acq On : 12 Aug 2022 13:22
 Operator : CG/JU
 Sample : PB146950BS
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS950

Quant Time: Aug 12 17:32:19 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	7.801	152	2087	0.400	ng/ul	0.02
4) Naphthalene-d8	10.591	136	7733	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.438	164	4530	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.197	188	8030	0.400	ng/ul	0.00
17) Chrysene-d12	21.401	240	5368	0.400	ng/ul	0.00
23) Perylene-d12	23.730	264	4115	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	1934	0.724	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.197	152	3668	0.324	ng/ul	0.02
18) Fluoranthene-d10	19.231	212	7564	0.456	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.232	88	2578	0.937	ng/ul	97
5) Naphthalene	10.640	128	3493	0.147	ng/ul	95
7) 2-Methylnaphthalene	12.268	142	2025	0.141	ng/ul	99
8) 1-Methylnaphthalene	12.482	142	2289	0.159	ng/ul	100
10) Acenaphthylene	14.165	152	3055	0.144	ng/ul	98
11) Acenaphthene	14.503	153	2525	0.144	ng/ul	99
12) Fluorene	15.502	166	2600	0.136	ng/ul	99
14) Pentachlorophenol	16.872	266	182	0.147	ng/ul	95
15) Phenanthrene	17.239	178	4104	0.143	ng/ul	98
16) Anthracene	17.336	178	3116	0.140	ng/ul	95
19) Fluoranthene	19.263	202	4511	0.190	ng/ul	95
20) Pyrene	19.626	202	4762	0.194	ng/ul	96
21) Benzo(a)anthracene	21.389	228	2459	0.140	ng/ul	97
22) Chrysene	21.439	228	3447	0.149	ng/ul	98
24) Benzo(b)fluoranthene	23.026	252	2821	0.148	ng/ul	81
25) Benzo(k)fluoranthene	23.072	252	3037	0.151	ng/ul#	83
26) Benzo(a)pyrene	23.634	252	2726	0.157	ng/ul#	66
27) Indeno(1,2,3-cd)pyrene	26.152	276	2833	0.137	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.172	278	2169	0.132	ng/ul#	79
29) Benzo(g,h,i)perylene	26.883	276	2756	0.149	ng/ul#	89

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

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