

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081222\
 Data File : BN021184.D
 Acq On : 12 Aug 2022 20:56
 Operator : CG/JU
 Sample : PB146950BS
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS950

Quant Time: Aug 13 01:03:15 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	2596	0.400	ng/ul	0.01
4) Naphthalene-d8	10.590	136	9246	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.437	164	5426	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.192	188	9813	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	6851	0.400	ng/ul	0.00
23) Perylene-d12	23.722	264	5349	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2230	0.671	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	4554	0.336	ng/ul	0.01
18) Fluoranthene-d10	19.229	212	9273	0.438	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	2908	0.850	ng/ul#	94
5) Naphthalene	10.639	128	4208	0.148	ng/ul	97
7) 2-Methylnaphthalene	12.267	142	2462	0.143	ng/ul	99
8) 1-Methylnaphthalene	12.481	142	2798	0.162	ng/ul	100
10) Acenaphthylene	14.159	152	3711	0.146	ng/ul	97
11) Acenaphthene	14.502	153	3032	0.144	ng/ul	99
12) Fluorene	15.501	166	3163	0.138	ng/ul	98
14) Pentachlorophenol	16.866	266	224	0.148	ng/ul	96
15) Phenanthrene	17.234	178	5017	0.143	ng/ul	99
16) Anthracene	17.335	178	3914	0.144	ng/ul	96
19) Fluoranthene	19.262	202	5538	0.183	ng/ul	99
20) Pyrene	19.625	202	5785	0.185	ng/ul	96
21) Benzo(a)anthracene	21.381	228	3335	0.149	ng/ul	97
22) Chrysene	21.430	228	4332	0.147	ng/ul	98
24) Benzo(b)fluoranthene	23.020	252	3623	0.146	ng/ul	82
25) Benzo(k)fluoranthene	23.067	252	3938	0.151	ng/ul#	88
26) Benzo(a)pyrene	23.626	252	3499	0.155	ng/ul#	66
27) Indeno(1,2,3-cd)pyrene	26.139	276	3686	0.137	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.159	278	2806	0.132	ng/ul#	84
29) Benzo(g,h,i)perylene	26.877	276	3445	0.143	ng/ul	93

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

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