

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021007.D
 Acq On : 05 Aug 2022 02:18
 Operator : CG/JU
 Sample : PB146744BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS744

Quant Time: Aug 05 03:25:27 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 : Wed Aug 03 16:31:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	2450	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7833	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	4431	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	8535	0.400	ng/ul	0.00
17) Chrysene-d12	21.415	240	7057	0.400	ng/ul	0.00
23) Perylene-d12	23.754	264	5824	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2040	0.651	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4076	0.355	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	8198	0.376	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	5474	1.695	ng/ul#	83
5) Naphthalene	10.651	128	7535	0.313	ng/ul	100
7) 2-Methylnaphthalene	12.279	142	4514	0.309	ng/ul	100
8) 1-Methylnaphthalene	12.494	142	4971	0.340	ng/ul	100
10) Acenaphthylene	14.174	152	6416	0.310	ng/ul	99
11) Acenaphthene	14.517	153	5178	0.302	ng/ul	98
12) Fluorene	15.511	166	5682	0.303	ng/ul	99
14) Pentachlorophenol	16.876	266	876	0.666	ng/ul	99
15) Phenanthrene	17.252	178	9043	0.296	ng/ul	100
16) Anthracene	17.345	178	7366	0.311	ng/ul	100
19) Fluoranthene	19.277	202	9982	0.320	ng/ul	99
20) Pyrene	19.640	202	10263	0.318	ng/ul	97
21) Benzo(a)anthracene	21.398	228	7581	0.329	ng/ul	99
22) Chrysene	21.450	228	9464	0.312	ng/ul	100
24) Benzo(b)fluoranthene	23.046	252	8631	0.320	ng/ul	99
25) Benzo(k)fluoranthene	23.090	252	8689	0.305	ng/ul	99
26) Benzo(a)pyrene	23.651	252	7809	0.318	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.176	276	8334	0.284	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.196	278	6635	0.286	ng/ul	99
29) Benzo(g,h,i)perylene	26.917	276	7580	0.289	ng/ul	99

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

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