

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021023.D
 Acq On : 05 Aug 2022 12:47
 Operator : CG/JU
 Sample : PB146649BS
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS649

Quant Time: Aug 06 00:27:51 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	2284	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7439	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.452	164	4165	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.205	188	8005	0.400	ng/ul	0.00
17) Chrysene-d12	21.415	240	6657	0.400	ng/ul	0.00
23) Perylene-d12	23.757	264	5498	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.207	96	2323	0.795	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4889	0.449	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	10165	0.494	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	5928	1.969	ng/ul#	84
5) Naphthalene	10.651	128	9612	0.421	ng/ul	99
7) 2-Methylnaphthalene	12.274	142	5827	0.420	ng/ul	99
8) 1-Methylnaphthalene	12.494	142	5991	0.432	ng/ul	99
10) Acenaphthylene	14.174	152	8217	0.422	ng/ul	99
11) Acenaphthene	14.517	153	6672	0.414	ng/ul	99
12) Fluorene	15.511	166	7381	0.419	ng/ul	99
14) Pentachlorophenol	16.872	266	1111	0.900	ng/ul	99
15) Phenanthrene	17.248	178	11943	0.417	ng/ul	100
16) Anthracene	17.345	178	9834	0.443	ng/ul	99
19) Fluoranthene	19.277	202	13167	0.447	ng/ul	99
20) Pyrene	19.640	202	13535	0.445	ng/ul	98
21) Benzo(a)anthracene	21.401	228	9736	0.448	ng/ul	100
22) Chrysene	21.453	228	12305	0.430	ng/ul	99
24) Benzo(b)fluoranthene	23.046	252	10917	0.429	ng/ul	99
25) Benzo(k)fluoranthene	23.093	252	11539	0.430	ng/ul	98
26) Benzo(a)pyrene	23.654	252	9972	0.429	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.179	276	10698	0.386	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.192	278	8548	0.390	ng/ul	97
29) Benzo(g,h,i)perylene	26.917	276	9873	0.399	ng/ul	100

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

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