

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020967.D
 Acq On : 03 Aug 2022 06:38
 Operator : CG/JU
 Sample : PB146670BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS670

Quant Time: Aug 03 07:08:48 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 04:19:04 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	2416	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7960	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	4391	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	8574	0.400	ng/ul	0.00
17) Chrysene-d12	21.415	240	7367	0.400	ng/ul	0.00
23) Perylene-d12	23.757	264	6522	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2105	0.714	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.208	152	4327	0.369	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	9013	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	5619	1.771	ng/ul#	86
5) Naphthalene	10.651	128	7971	0.366	ng/ul	100
7) 2-Methylnaphthalene	12.279	142	4726	0.357	ng/ul	100
8) 1-Methylnaphthalene	12.494	142	5134	0.351	ng/ul	99
10) Acenaphthylene	14.179	152	6670	0.372	ng/ul	99
11) Acenaphthene	14.521	153	5493	0.366	ng/ul	98
12) Fluorene	15.511	166	6084	0.352	ng/ul	100
14) Pentachlorophenol	16.876	266	924	0.675	ng/ul	98
15) Phenanthrene	17.252	178	9846	0.339	ng/ul	100
16) Anthracene	17.345	178	7968	0.319	ng/ul	100
19) Fluoranthene	19.277	202	10879	0.329	ng/ul	100
20) Pyrene	19.640	202	11168	0.322	ng/ul	98
21) Benzo(a)anthracene	21.401	228	8463	0.316	ng/ul	100
22) Chrysene	21.453	228	10666	0.322	ng/ul	100
24) Benzo(b)fluoranthene	23.046	252	10219	0.328	ng/ul	99
25) Benzo(k)fluoranthene	23.093	252	10498	0.311	ng/ul	99
26) Benzo(a)pyrene	23.654	252	9429	0.321	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.179	276	10475	0.303	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.196	278	8397	0.297	ng/ul	97
29) Benzo(g,h,i)perylene	26.917	276	9167	0.292	ng/ul	100

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

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