

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021008.D
 Acq On : 05 Aug 2022 02:55
 Operator : CG/JU
 Sample : PB146574BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS574

Quant Time: Aug 05 03:25:42 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	7.809	152	2315	0.400	ng/ul	0.00
4) Naphthalene-d8	10.601	136	7412	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.456	164	4172	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.209	188	8031	0.400	ng/ul	0.00
17) Chrysene-d12	21.412	240	6587	0.400	ng/ul	0.00
23) Perylene-d12	23.756	264	5586	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	2084	0.704	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4257	0.392	ng/ul	0.00
18) Fluoranthene-d10	19.244	212	8615	0.424	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	5674	1.860	ng/ul#	84
5) Naphthalene	10.651	128	7605	0.334	ng/ul	100
7) 2-Methylnaphthalene	12.279	142	4636	0.336	ng/ul	100
8) 1-Methylnaphthalene	12.493	142	5083	0.368	ng/ul	100
10) Acenaphthylene	14.174	152	6542	0.335	ng/ul	99
11) Acenaphthene	14.516	153	5374	0.333	ng/ul	99
12) Fluorene	15.511	166	5834	0.331	ng/ul	99
14) Pentachlorophenol	16.876	266	834	0.673	ng/ul	99
15) Phenanthrene	17.251	178	9411	0.328	ng/ul	100
16) Anthracene	17.349	178	7634	0.343	ng/ul	99
19) Fluoranthene	19.277	202	10334	0.354	ng/ul	99
20) Pyrene	19.639	202	10644	0.354	ng/ul	97
21) Benzo(a)anthracene	21.397	228	7861	0.366	ng/ul	99
22) Chrysene	21.450	228	9811	0.347	ng/ul	100
24) Benzo(b)fluoranthene	23.043	252	9147	0.354	ng/ul	98
25) Benzo(k)fluoranthene	23.090	252	9246	0.339	ng/ul	99
26) Benzo(a)pyrene	23.651	252	8290	0.351	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.175	276	9035	0.321	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.192	278	7166	0.322	ng/ul	98
29) Benzo(g,h,i)perylene	26.913	276	8050	0.320	ng/ul	99

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

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