

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082222\
 Data File : BN021308.D
 Acq On : 22 Aug 2022 20:02
 Operator : CG/JU
 Sample : PB147154BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS154

Quant Time: Aug 23 04:03:33 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 18:06:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.097	152	2560	0.400	ng/ul	0.00
4) Naphthalene-d8	10.933	136	8210	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.754	164	4406	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.502	188	9265	0.400	ng/ul	0.00
17) Chrysene-d12	21.688	240	7595	0.400	ng/ul	# 0.00
23) Perylene-d12	24.222	264	6200	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.384	96	2052	0.621	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.517	152	3958	0.319	ng/ul	0.00
18) Fluoranthene-d10	19.524	212	7527	0.322	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.422	88	5560	1.607	ng/ul#	88
5) Naphthalene	10.982	128	7409	0.285	ng/ul	98
7) 2-Methylnaphthalene	12.588	142	4661	0.287	ng/ul	98
8) 1-Methylnaphthalene	12.808	142	4879	0.320	ng/ul	99
10) Acenaphthylene	14.471	152	6191	0.272	ng/ul	99
11) Acenaphthene	14.814	153	4829	0.277	ng/ul	98
12) Fluorene	15.799	166	5440	0.280	ng/ul	100
14) Pentachlorophenol	17.143	266	477	0.466	ng/ul	98
15) Phenanthrene	17.544	178	9421	0.274	ng/ul	99
16) Anthracene	17.633	178	7410	0.269	ng/ul	100
19) Fluoranthene	19.557	202	10116	0.281	ng/ul	99
20) Pyrene	19.919	202	10109	0.281	ng/ul	97
21) Benzo(a)anthracene	21.673	228	7721	0.293	ng/ul	99
22) Chrysene	21.726	228	9988	0.295	ng/ul	99
24) Benzo(b)fluoranthene	23.445	252	8657	0.325	ng/ul	97
25) Benzo(k)fluoranthene	23.494	252	9099	0.317	ng/ul#	96
26) Benzo(a)pyrene	24.108	252	7703	0.307	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.894	276	9467	0.321	ng/ul#	42
28) Dibenzo(a,h)anthracene	26.924	278	7753	0.320	ng/ul	95
29) Benzo(g,h,i)perylene	27.723	276	9233	0.316	ng/ul	98

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

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