

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN063022\
 Data File : BN020591.D
 Acq On : 30 Jun 2022 13:53
 Operator : CG/JU
 Sample : PB145665BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS665

Quant Time: Jun 30 15:41:28 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	3399	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.524	136	10318	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.368	164	5416	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.111	188	10544	0.400	ng/ul	0.00
17) Chrysene-d12	21.305	240	7514	0.400	ng/ul	0.00
23) Perylene-d12	23.585	264	5553	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.197	96	3177	0.807	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.119	152	6079	0.359	ng/ul	0.00
18) Fluoranthene-d10	19.141	212	11388	0.460	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.227	88	8257	2.054	ng/ul#	79
5) Naphthalene	10.568	128	11657	0.366	ng/ul	99
7) 2-Methylnaphthalene	12.190	142	7031	0.330	ng/ul	99
8) 1-Methylnaphthalene	12.410	142	7426	0.367	ng/ul	99
10) Acenaphthylene	14.090	152	9276	0.361	ng/ul	100
11) Acenaphthene	14.432	153	7528	0.366	ng/ul	99
12) Fluorene	15.418	166	8325	0.342	ng/ul	99
14) Pentachlorophenol	16.782	266	833	0.392	ng/ul	98
15) Phenanthrene	17.154	178	13740	0.375	ng/ul	100
16) Anthracene	17.247	178	11180	0.353	ng/ul	99
19) Fluoranthene	19.174	202	14651	0.427	ng/ul	99
20) Pyrene	19.536	202	14909	0.419	ng/ul	100
21) Benzo(a)anthracene	21.290	228	9383	0.348	ng/ul	99
22) Chrysene	21.340	228	12132	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.898	252	10039	0.391	ng/ul	98
25) Benzo(k)fluoranthene	22.945	252	10615	0.409	ng/ul	99
26) Benzo(a)pyrene	23.491	252	9199	0.394	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.918	276	9768	0.420	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.931	278	7788	0.415	ng/ul	98
29) Benzo(g,h,i)perylene	26.632	276	8990	0.451	ng/ul	99

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

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