

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071922\
 Data File : BN020792.D
 Acq On : 20 Jul 2022 02:28
 Operator : CG/JU
 Sample : PB146297BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS297

Quant Time: Jul 20 03:13:49 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 : Wed Jul 20 01:02:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	2162	0.400	ng/ul	0.00
4) Naphthalene-d8	10.617	136	8405	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.469	164	4691	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	9503	0.400	ng/ul	0.00
17) Chrysene-d12	21.422	240	10680	0.400	ng/ul	0.00
23) Perylene-d12	23.772	264	8698	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	2291	0.915	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	5913	0.429	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	11889	0.338	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	5603	2.191	ng/ul#	82
5) Naphthalene	10.667	128	10606	0.409	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	6817	0.392	ng/ul	99
8) 1-Methylnaphthalene	12.509	142	7264	0.441	ng/ul	100
10) Acenaphthylene	14.192	152	10089	0.454	ng/ul	99
11) Acenaphthene	14.534	153	7163	0.402	ng/ul	99
12) Fluorene	15.524	166	8216	0.389	ng/ul	99
14) Pentachlorophenol	16.888	266	1372	0.716	ng/ul	97
15) Phenanthrene	17.263	178	13956	0.422	ng/ul	100
16) Anthracene	17.356	178	11073	0.388	ng/ul	100
19) Fluoranthene	19.290	202	14995	0.307	ng/ul	99
20) Pyrene	19.653	202	15626	0.309	ng/ul	99
21) Benzo(a)anthracene	21.407	228	15017	0.391	ng/ul	99
22) Chrysene	21.460	228	19164	0.449	ng/ul	100
24) Benzo(b)fluoranthene	23.053	252	17700	0.440	ng/ul	97
25) Benzo(k)fluoranthene	23.102	252	18858	0.464	ng/ul	99
26) Benzo(a)pyrene	23.664	252	16025	0.438	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.200	276	17495	0.481	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.216	278	13743	0.467	ng/ul	99
29) Benzo(g,h,i)perylene	26.941	276	15928	0.510	ng/ul	100

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

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