

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071922\
 Data File : BN020781.D
 Acq On : 19 Jul 2022 19:01
 Operator : CG/JU
 Sample : PB146235BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS235

Quant Time: Jul 20 01:05:33 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	7.830	152	1870	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	7449	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	3779	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.226	188	7710	0.400	ng/ul	0.00
17) Chrysene-d12	21.432	240	9171	0.400	ng/ul	0.00
23) Perylene-d12	23.779	264	7462	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	1922	0.887	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	4791	0.392	ng/ul	0.00
18) Fluoranthene-d10	19.262	212	12167	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	5062	2.288	ng/ul#	82
5) Naphthalene	10.672	128	9449	0.411	ng/ul	100
7) 2-Methylnaphthalene	12.295	142	5970	0.388	ng/ul	99
8) 1-Methylnaphthalene	12.515	142	6399	0.439	ng/ul	99
10) Acenaphthylene	14.192	152	7849	0.438	ng/ul	100
11) Acenaphthene	14.534	153	6053	0.422	ng/ul	100
12) Fluorene	15.529	166	6993	0.412	ng/ul	99
14) Pentachlorophenol	16.892	266	947	0.609	ng/ul	99
15) Phenanthrene	17.268	178	11141	0.416	ng/ul	100
16) Anthracene	17.365	178	9260	0.400	ng/ul	100
19) Fluoranthene	19.295	202	16936	0.404	ng/ul	100
20) Pyrene	19.653	202	17492	0.403	ng/ul	97
21) Benzo(a)anthracene	21.417	228	12873	0.391	ng/ul	100
22) Chrysene	21.467	228	16719	0.456	ng/ul	98
24) Benzo(b)fluoranthene	23.066	252	15417	0.446	ng/ul	98
25) Benzo(k)fluoranthene	23.112	252	16644	0.477	ng/ul	99
26) Benzo(a)pyrene	23.677	252	13770	0.439	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.215	276	15431	0.494	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.238	278	11959	0.474	ng/ul	99
29) Benzo(g,h,i)perylene	26.959	276	13985	0.522	ng/ul	99

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

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