

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072623\
 Data File : BN026759.D
 Acq On : 26 Jul 2023 16:46
 Operator : MA/JU
 Sample : PB154053BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS053

Quant Time: Jul 26 20:21:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 00:39:11 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	1690	0.400	ng/ul	0.00
4) Naphthalene-d8	10.630	136	5180	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.462	164	3527	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.215	188	7516	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.422	240	7873	0.400	ng/ul	# 0.00
23) Perylene-d12	23.796	264	7694	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.191	96	1883	0.846	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.231	152	2569	0.365	ng/ul	0.01
18) Fluoranthene-d10	19.250	212	8737	0.320	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.224	88	4182	1.986	ng/ul#	88
5) Naphthalene	10.680	128	4836	0.340	ng/ul	96
7) 2-Methylnaphthalene	12.302	142	2781	0.345	ng/ul	99
8) 1-Methylnaphthalene	12.511	142	3315	0.383	ng/ul	99
10) Acenaphthylene	14.185	152	5871	0.340	ng/ul	99
11) Acenaphthene	14.522	153	4291	0.326	ng/ul	99
12) Fluorene	15.517	166	4834	0.350	ng/ul	98
14) Pentachlorophenol	16.894	266	1176	0.620	ng/ul	98
15) Phenanthrene	17.257	178	7796	0.326	ng/ul	98
16) Anthracene	17.358	178	7326	0.353	ng/ul	99
19) Fluoranthene	19.278	202	10738	0.291	ng/ul#	92
20) Pyrene	19.641	202	11603	0.290	ng/ul#	92
21) Benzo(a)anthracene	21.405	228	9397	0.325	ng/ul	99
22) Chrysene	21.454	228	11017	0.335	ng/ul	99
24) Benzo(b)fluoranthene	23.074	252	10509	0.351	ng/ul	93
25) Benzo(k)fluoranthene	23.123	252	11792	0.375	ng/ul#	95
26) Benzo(a)pyrene	23.690	252	9188	0.344	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.260	276	10265	0.293	ng/ul#	72
28) Dibenzo(a,h)anthracene	26.271	278	7688	0.286	ng/ul#	89
29) Benzo(g,h,i)perylene	27.005	276	9342	0.303	ng/ul#	88

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

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