

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021723\
 Data File : BN024064.D
 Acq On : 17 Feb 2023 13:24
 Operator : CG/JU
 Sample : PB150872BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS872

Quant Time: Feb 17 21:35:46 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 14:45:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	4840	0.400	ng/ul	0.00
4) Naphthalene-d8	10.634	136	12665	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.483	164	7297	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.234	188	15854	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.422	240	13545	0.400	ng/ul	# 0.00
23) Perylene-d12	23.804	264	9836	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.226	96	2541	0.490	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.228	152	4939	0.249	ng/ul	0.00
18) Fluoranthene-d10	19.262	212	11808	0.325	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.260	88	7379	1.342	ng/ul#	80
5) Naphthalene	10.683	128	9427	0.272	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	5903	0.262	ng/ul	100
8) 1-Methylnaphthalene	12.525	142	6198	0.269	ng/ul	100
10) Acenaphthylene	14.201	152	7559	0.271	ng/ul	99
11) Acenaphthene	14.543	153	6778	0.276	ng/ul	98
12) Fluorene	15.533	166	7702	0.259	ng/ul	97
14) Pentachlorophenol	16.888	266	1929	0.369	ng/ul	99
15) Phenanthrene	17.276	178	13109	0.277	ng/ul	100
16) Anthracene	17.365	178	10850	0.274	ng/ul	98
19) Fluoranthene	19.295	202	15675	0.348	ng/ul	96
20) Pyrene	19.657	202	15824	0.339	ng/ul#	94
21) Benzo(a)anthracene	21.407	228	13489	0.293	ng/ul	99
22) Chrysene	21.457	228	15136	0.312	ng/ul	99
24) Benzo(b)fluoranthene	23.079	252	14254	0.328	ng/ul	96
25) Benzo(k)fluoranthene	23.126	252	13906	0.327	ng/ul#	95
26) Benzo(a)pyrene	23.696	252	10991	0.304	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.250	276	13323	0.262	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.270	278	10906	0.269	ng/ul#	95
29) Benzo(g,h,i)perylene	27.011	276	11874	0.284	ng/ul	97

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

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