

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029674.D
 Acq On : 11 Feb 2024 08:04
 Operator : MA/JU
 Sample : PB158928BS
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS928

Quant Time: Feb 11 23:25:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Feb 11 23:18:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.880	152	5493	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	15173	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	6475	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	11412	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	8268	0.400	ng/ul	0.00
23) Perylene-d12	23.843	264	8256	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	5303	0.774	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.273	152	7676	0.398	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	10816	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.366	88	17025	2.283	ng/ul#	88
5) Naphthalene	10.727	128	17112	0.401	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	9633	0.393	ng/ul	99
8) 1-Methylnaphthalene	12.564	142	9555	0.383	ng/ul	100
10) Acenaphthylene	14.243	152	12511	0.396	ng/ul	99
11) Acenaphthene	14.585	153	9624	0.401	ng/ul	100
12) Fluorene	15.575	166	9865	0.393	ng/ul	99
14) Pentachlorophenol	16.917	266	2049	0.805	ng/ul#	50
15) Phenanthrene	17.314	178	14218	0.399	ng/ul	99
16) Anthracene	17.407	178	12074	0.403	ng/ul	99
19) Fluoranthene	19.337	202	14768	0.389	ng/ul	99
20) Pyrene	19.699	202	15251	0.392	ng/ul	99
21) Benzo(a)anthracene	21.455	228	12224	0.410	ng/ul	99
22) Chrysene	21.508	228	13583	0.416	ng/ul	99
24) Benzo(b)fluoranthene	23.118	252	13537	0.436	ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	15078	0.438	ng/ul	99
26) Benzo(a)pyrene	23.735	252	12381	0.424	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.292	276	15529	0.438	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.325	278	12230	0.440	ng/ul	99
29) Benzo(g,h,i)perylene	27.040	276	14236	0.442	ng/ul	98

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

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