

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029692.D
 Acq On : 11 Feb 2024 19:32
 Operator : MA/JU
 Sample : PB158926BS
 Misc :
 ALS Vial : 85 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS926

Quant Time: Feb 11 23:28:05 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	4988	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	13718	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	5942	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	10979	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	7958	0.400	ng/ul	0.00
23) Perylene-d12	23.840	264	7905	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	5299	0.852	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	7352	0.421	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	11175	0.427	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.361	88	16741	2.472	ng/ul#	85
5) Naphthalene	10.727	128	16436	0.426	ng/ul	100
7) 2-Methylnaphthalene	12.344	142	9381	0.424	ng/ul	100
8) 1-Methylnaphthalene	12.558	142	9170	0.407	ng/ul	99
10) Acenaphthylene	14.238	152	12351	0.426	ng/ul	99
11) Acenaphthene	14.580	153	9492	0.431	ng/ul	98
12) Fluorene	15.570	166	9918	0.430	ng/ul	99
14) Pentachlorophenol	16.917	266	2201	0.899	ng/ul#	51
15) Phenanthrene	17.314	178	14758	0.430	ng/ul	100
16) Anthracene	17.407	178	12500	0.433	ng/ul	100
19) Fluoranthene	19.332	202	15364	0.420	ng/ul	99
20) Pyrene	19.699	202	15872	0.424	ng/ul	98
21) Benzo(a)anthracene	21.455	228	12445	0.434	ng/ul	99
22) Chrysene	21.508	228	13937	0.444	ng/ul	100
24) Benzo(b)fluoranthene	23.118	252	13741	0.462	ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	15060	0.457	ng/ul	98
26) Benzo(a)pyrene	23.735	252	12428	0.444	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.292	276	15591	0.459	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.322	278	12190	0.458	ng/ul	98
29) Benzo(g,h,i)perylene	27.036	276	14464	0.469	ng/ul	99

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

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