

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029703.D
 Acq On : 12 Feb 2024 02:47
 Operator : MA/JU
 Sample : PB158924BS
 Misc :
 ALS Vial : 97 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS924

Quant Time: Feb 12 03:17:49 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	5018	0.400	ng/ul	0.00
4) Naphthalene-d8	10.672	136	13559	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	5728	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	9931	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	6998	0.400	ng/ul	0.00
23) Perylene-d12	23.840	264	6922	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	5307	0.848	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	7439	0.431	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	10338	0.450	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.366	88	16761	2.461	ng/ul#	88
5) Naphthalene	10.722	128	16654	0.436	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	9411	0.430	ng/ul	100
8) 1-Methylnaphthalene	12.559	142	9224	0.414	ng/ul	99
10) Acenaphthylene	14.238	152	12016	0.430	ng/ul	98
11) Acenaphthene	14.581	153	9325	0.439	ng/ul	98
12) Fluorene	15.571	166	9423	0.424	ng/ul	99
14) Pentachlorophenol	16.917	266	1976	0.892	ng/ul#	51
15) Phenanthrene	17.314	178	13668	0.441	ng/ul	100
16) Anthracene	17.407	178	11505	0.441	ng/ul	99
19) Fluoranthene	19.332	202	14089	0.438	ng/ul	99
20) Pyrene	19.699	202	14566	0.442	ng/ul	99
21) Benzo(a)anthracene	21.452	228	11265	0.446	ng/ul	100
22) Chrysene	21.505	228	12883	0.467	ng/ul	99
24) Benzo(b)fluoranthene	23.118	252	12597	0.484	ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	13981	0.484	ng/ul	97
26) Benzo(a)pyrene	23.732	252	11639	0.475	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.289	276	14458	0.487	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.319	278	11194	0.480	ng/ul	98
29) Benzo(g,h,i)perylene	27.037	276	13486	0.500	ng/ul	99

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

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