

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029702.D
 Acq On : 12 Feb 2024 02:10
 Operator : MA/JU
 Sample : PB158920BS
 Misc :
 ALS Vial : 96 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS920

Quant Time: Feb 12 02:42:13 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	5199	0.400	ng/ul	0.00
4) Naphthalene-d8	10.672	136	14022	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	6047	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	10885	0.400	ng/ul	0.00
17) Chrysene-d12	21.469	240	7935	0.400	ng/ul	0.00
23) Perylene-d12	23.839	264	7828	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	4933	0.761	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	7009	0.393	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	10230	0.392	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.366	88	15494	2.195	ng/ul#	87
5) Naphthalene	10.727	128	15529	0.394	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	8720	0.385	ng/ul	99
8) 1-Methylnaphthalene	12.559	142	8644	0.375	ng/ul	99
10) Acenaphthylene	14.238	152	11331	0.384	ng/ul	99
11) Acenaphthene	14.581	153	8761	0.391	ng/ul	98
12) Fluorene	15.571	166	8982	0.383	ng/ul	99
14) Pentachlorophenol	16.917	266	1982	0.816	ng/ul#	50
15) Phenanthrene	17.314	178	13181	0.388	ng/ul	99
16) Anthracene	17.407	178	11085	0.387	ng/ul	99
19) Fluoranthene	19.332	202	13977	0.383	ng/ul	99
20) Pyrene	19.694	202	14438	0.387	ng/ul	98
21) Benzo(a)anthracene	21.454	228	11488	0.402	ng/ul	99
22) Chrysene	21.507	228	13084	0.418	ng/ul	100
24) Benzo(b)fluoranthene	23.117	252	12968	0.440	ng/ul	98
25) Benzo(k)fluoranthene	23.167	252	14107	0.432	ng/ul	98
26) Benzo(a)pyrene	23.734	252	11721	0.423	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.290	276	14495	0.431	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.320	278	11217	0.425	ng/ul	96
29) Benzo(g,h,i)perylene	27.031	276	13418	0.440	ng/ul	99

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

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