

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021324\
 Data File : BN029719.D
 Acq On : 12 Feb 2024 18:26
 Operator : MA/JU
 Sample : PB158957BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS957

Quant Time: Feb 12 23:34:35 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	7.876	152	5002	0.400	ng/ul	0.00
4) Naphthalene-d8	10.672	136	13401	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.516	164	5630	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	9756	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	7000	0.400	ng/ul	0.00
23) Perylene-d12	23.840	264	6712	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	4630	0.742	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	6447	0.378	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	8845	0.384	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.361	88	14655	2.158	ng/ul#	87
5) Naphthalene	10.722	128	14289	0.379	ng/ul	99
7) 2-Methylnaphthalene	12.338	142	7995	0.370	ng/ul	100
8) 1-Methylnaphthalene	12.558	142	7919	0.360	ng/ul	99
10) Acenaphthylene	14.238	152	10326	0.376	ng/ul	99
11) Acenaphthene	14.580	153	7902	0.378	ng/ul	99
12) Fluorene	15.570	166	7978	0.365	ng/ul	99
14) Pentachlorophenol	16.917	266	1659	0.762	ng/ul#	50
15) Phenanthrene	17.314	178	11642	0.382	ng/ul	99
16) Anthracene	17.407	178	9774	0.381	ng/ul	99
19) Fluoranthene	19.332	202	12061	0.375	ng/ul	98
20) Pyrene	19.699	202	12480	0.379	ng/ul	98
21) Benzo(a)anthracene	21.455	228	9432	0.374	ng/ul	99
22) Chrysene	21.508	228	11079	0.401	ng/ul	99
24) Benzo(b)fluoranthene	23.121	252	10892	0.431	ng/ul	99
25) Benzo(k)fluoranthene	23.171	252	11661	0.416	ng/ul	97
26) Benzo(a)pyrene	23.735	252	9724	0.409	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.292	276	12214	0.424	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.325	278	9438	0.417	ng/ul	95
29) Benzo(g,h,i)perylene	27.040	276	11527	0.441	ng/ul	99

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

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