

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029675.D
 Acq On : 11 Feb 2024 08:41
 Operator : MA/JU
 Sample : PB158930BS
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS930

Quant Time: Feb 11 23:25:40 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	5270	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	14311	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	6150	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	11178	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	8113	0.400	ng/ul	0.00
23) Perylene-d12	23.840	264	7933	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	5042	0.767	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.272	152	7084	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	10384	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.365	88	16207	2.266	ng/ul#	88
5) Naphthalene	10.727	128	15869	0.394	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	8952	0.388	ng/ul	99
8) 1-Methylnaphthalene	12.564	142	8833	0.376	ng/ul	100
10) Acenaphthylene	14.243	152	11706	0.390	ng/ul	99
11) Acenaphthene	14.585	153	9006	0.395	ng/ul	99
12) Fluorene	15.575	166	9300	0.390	ng/ul	98
14) Pentachlorophenol	16.917	266	2025	0.812	ng/ul#	50
15) Phenanthrene	17.314	178	13681	0.392	ng/ul	100
16) Anthracene	17.407	178	11606	0.395	ng/ul	99
19) Fluoranthene	19.336	202	14277	0.383	ng/ul	99
20) Pyrene	19.699	202	14643	0.383	ng/ul	99
21) Benzo(a)anthracene	21.455	228	11778	0.403	ng/ul	99
22) Chrysene	21.508	228	13173	0.411	ng/ul	100
24) Benzo(b)fluoranthene	23.121	252	12966	0.434	ng/ul	100
25) Benzo(k)fluoranthene	23.168	252	14504	0.438	ng/ul	98
26) Benzo(a)pyrene	23.738	252	11695	0.417	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.295	276	14961	0.439	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.322	278	11706	0.438	ng/ul	97
29) Benzo(g,h,i)perylene	27.043	276	13744	0.444	ng/ul	98

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

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