

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030131.D
 Acq On : 28 Feb 2024 15:25
 Operator : MA/JU
 Sample : PB159157BS
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS157

Quant Time: Feb 28 15:58:24 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 : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	2867	0.400	ng/ul	0.00
4) Naphthalene-d8	10.650	136	8157	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	3968	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.259	188	6825	0.400	ng/ul	0.00
17) Chrysene-d12	21.461	240	4912	0.400	ng/ul	0.00
23) Perylene-d12	23.820	264	4666	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	2776	0.777	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	3993	0.385	ng/ul	0.00
18) Fluoranthene-d10	19.290	212	6681	0.414	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.349	88	7536	1.936	ng/ul#	86
5) Naphthalene	10.700	128	8150	0.355	ng/ul	99
7) 2-Methylnaphthalene	12.322	142	4771	0.362	ng/ul	99
8) 1-Methylnaphthalene	12.537	142	5034	0.376	ng/ul	99
10) Acenaphthylene	14.220	152	6996	0.361	ng/ul	98
11) Acenaphthene	14.562	153	5412	0.368	ng/ul	98
12) Fluorene	15.557	166	5423	0.352	ng/ul	96
14) Pentachlorophenol	16.909	266	1403	0.922	ng/ul#	50
15) Phenanthrene	17.302	178	7687	0.361	ng/ul	96
16) Anthracene	17.394	178	6757	0.377	ng/ul	97
19) Fluoranthene	19.318	202	8878	0.393	ng/ul	97
20) Pyrene	19.680	202	9478	0.410	ng/ul	94
21) Benzo(a)anthracene	21.444	228	6374	0.360	ng/ul	98
22) Chrysene	21.496	228	7543	0.389	ng/ul	97
24) Benzo(b)fluoranthene	23.101	252	6910	0.394	ng/ul	95
25) Benzo(k)fluoranthene	23.150	252	8025	0.412	ng/ul	93
26) Benzo(a)pyrene	23.718	252	6633	0.402	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.279	276	8156	0.407	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.322	278	6348	0.404	ng/ul#	87
29) Benzo(g,h,i)perylene	27.010	276	7435	0.409	ng/ul	93

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

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