

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030135.D
 Acq On : 28 Feb 2024 18:25
 Operator : MA/JU
 Sample : PB159159BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS159

Quant Time: Feb 28 23:06:25 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	4753	0.400	ng/ul	0.00
4) Naphthalene-d8	10.650	136	13268	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	5756	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.259	188	10576	0.400	ng/ul	0.00
17) Chrysene-d12	21.458	240	8058	0.400	ng/ul	0.00
23) Perylene-d12	23.826	264	7128	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	4112	0.694	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	6002	0.356	ng/ul	0.00
18) Fluoranthene-d10	19.290	212	10410	0.393	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.353	88	11796	1.828	ng/ul#	85
5) Naphthalene	10.700	128	13212	0.354	ng/ul	99
7) 2-Methylnaphthalene	12.322	142	7273	0.340	ng/ul	99
8) 1-Methylnaphthalene	12.542	142	7760	0.356	ng/ul	99
10) Acenaphthylene	14.224	152	9740	0.347	ng/ul	98
11) Acenaphthene	14.562	153	7889	0.370	ng/ul	97
12) Fluorene	15.561	166	7941	0.356	ng/ul	98
14) Pentachlorophenol	16.909	266	1713	0.726	ng/ul#	48
15) Phenanthrene	17.301	178	11907	0.360	ng/ul	98
16) Anthracene	17.394	178	10053	0.362	ng/ul	98
19) Fluoranthene	19.323	202	13846	0.374	ng/ul	98
20) Pyrene	19.685	202	14707	0.388	ng/ul	97
21) Benzo(a)anthracene	21.443	228	10000	0.344	ng/ul	99
22) Chrysene	21.496	228	12509	0.393	ng/ul	100
24) Benzo(b)fluoranthene	23.106	252	11318	0.422	ng/ul	99
25) Benzo(k)fluoranthene	23.156	252	13410	0.451	ng/ul	97
26) Benzo(a)pyrene	23.723	252	10328	0.410	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.285	276	12090	0.395	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.332	278	9242	0.385	ng/ul#	92
29) Benzo(g,h,i)perylene	27.023	276	11313	0.407	ng/ul	96

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

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