

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070524\
 Data File : BN032510.D
 Acq On : 05 Jul 2024 15:15
 Operator : MA/JU
 Sample : PB161671BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS671

Quant Time: Jul 05 15:40:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.595	152	4241	0.400	ng/ul	0.00
4) Naphthalene-d8	10.371	136	12218	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.244	164	6566	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.007	188	11921	0.400	ng/ul	0.00
17) Chrysene-d12	21.217	240	11650	0.400	ng/ul	# 0.00
23) Perylene-d12	23.435	264	14293	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	4346	0.922	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.976	152	6342	0.346	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	12609	0.360	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.173	88	9196	1.714	ng/ul	90
5) Naphthalene	10.426	128	11445	0.354	ng/ul	97
7) 2-Methylnaphthalene	12.053	142	6680	0.312	ng/ul	100
8) 1-Methylnaphthalene	12.268	142	6632	0.298	ng/ul	99
10) Acenaphthylene	13.966	152	10798	0.395	ng/ul	97
11) Acenaphthene	14.308	153	7507	0.354	ng/ul	99
12) Fluorene	15.308	166	7868	0.312	ng/ul	99
14) Pentachlorophenol	16.682	266	1839	0.682	ng/ul	99
15) Phenanthrene	17.050	178	11593	0.355	ng/ul	98
16) Anthracene	17.147	178	11185	0.384	ng/ul	97
19) Fluoranthene	19.078	202	14622	0.330	ng/ul	97
20) Pyrene	19.440	202	15611	0.337	ng/ul	99
21) Benzo(a)anthracene	21.203	228	15757	0.375	ng/ul	97
22) Chrysene	21.252	228	17021	0.369	ng/ul	96
24) Benzo(b)fluoranthene	22.766	252	18920	0.305	ng/ul	87
25) Benzo(k)fluoranthene	22.810	252	18725	0.310	ng/ul#	87
26) Benzo(a)pyrene	23.336	252	17065	0.395	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.676	276	22210	0.328	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.680	278	17705	0.352	ng/ul#	86
29) Benzo(g,h,i)perylene	26.360	276	17857	0.331	ng/ul	92

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

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