

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100923\
 Data File : BM042245.D
 Acq On : 10 Oct 2023 04:00
 Operator : MA/JU
 Sample : PB156201BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS201

Quant Time: Oct 10 05:06:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 09 23:59:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.047	152	5	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.873	136	103	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.694	164	318	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.443	188	444	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.623	240	190	0.400	ng/ul	#-0.01
23) Perylene-d12	0.000	264	0	0.000	ng/ul	-24.14
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	43098	6472.675	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	4907	34.736	ng/ul	0.00
18) Fluoranthene-d10	19.464	212	7829	10.975	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.444	88	90291	12985.305	ng/ul#	73
5) Naphthalene	10.928	128	1326727	3860.982	ng/ul	98
7) 2-Methylnaphthalene	12.528	142	818532	3978.416	ng/ul	99
8) 1-Methylnaphthalene	12.748	142	782855	3777.372	ng/ul	99
10) Acenaphthylene	14.421	152	1382725	688.332	ng/ul	100
11) Acenaphthene	14.754	153	892297	586.810	ng/ul	99
12) Fluorene	15.739	166	990577	578.254	ng/ul	100
14) Pentachlorophenol	17.071	266	302609	1503.412	ng/ul	99
15) Phenanthrene	17.485	178	1532706	837.288	ng/ul	100
16) Anthracene	17.578	178	1508795	901.037	ng/ul	99
19) Fluoranthene	19.496	202	1714370	1187.825	ng/ul	97
20) Pyrene	19.863	202	1660315	1082.327	ng/ul	96
21) Benzo(a)anthracene	21.609	228	1295505	1056.190	ng/ul	99
22) Chrysene	21.664	228	1173526	982.970	ng/ul	99

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

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